



Analytical Resources, Incorporated
Analytical Chemists and Consultants



October 13, 2011

John Long
AMEC/Geomatrix
600 University Suite 600
Seattle, WA 98101

RE: Client Project: Former Rhone Poulenc- 8769 Shoreline Investigation
ARI Job Numbers: TN19, TN21

Dear John:

Please find enclosed the final data package for samples for the project referenced above. ARI received five water samples and one trip blank on September 20, 2011.

Please refer to the case narrative for details on the analyses of these samples.

An electronic copy of this package will be kept on file at ARI. If you have questions or problems, please feel free to contact me at any time.

Sincerely,

ANALYTICAL RESOURCES, INC.


Cheronne Oreiro
Project Manager
-For-
Kelly Bottem
Client Services Manager
206/695-6211
kellyb@arilabs.com

Enclosures

cc: file TN19_TN21



**Spike Recovery Control Limits for Analysis of Aqueous Samples
Volatile Organic Compounds (VOA) EPA SW-846 Methods 8260C
10 mL Purge Volume ^(1,6)**

Effective: 8/30/2010

Control limits are updated periodically. Assure that you have ARI's current control limits by downloading the files at the time of use. <http://www.arilabs.com/portal/downloads/ARI-CLS.zip>

LCS Spike Recovery ⁽⁵⁾	ARI Control Limits	ARI ME Control Limits ⁽²⁾
Dichlorodifluoromethane	69 - 122	60 - 131
Chloromethane	76 - 120	69 - 123
Vinyl Chloride	80 - 120	75 - 123
Bromomethane	71 - 120	63 - 129
Chloroethane	80 - 120	75 - 124
1,1,2-Trichloro-1,2,2-trifluoroethane	80 - 121	76 - 128
Acrolein	69 - 126	60 - 136
Acetone	71 - 120	64 - 120
1,1-Dichloroethene	80 - 120	79 - 122
Bromoethane	80 - 120	80 - 121
Methyl Iodide	76 - 120	69 - 127
Methylene Chloride	80 - 120	77 - 120
Acrylonitrile	79 - 120	74 - 120
Methyl tert-Butyl Ether	80 - 120	77 - 121
Carbon Disulfide	80 - 120	78 - 121
trans-1,2-Dichloroethene	80 - 120	80 - 120
Vinyl Acetate	80 - 120	76 - 120
1,1-Dichloroethane	80 - 120	80 - 120
2-Butanone	80 - 120	76 - 120
2,2-Dichloropropane	80 - 120	77 - 120
cis-1,2-Dichloroethene	80 - 120	80 - 120
Chloroform	80 - 120	80 - 120
Bromodichloromethane	80 - 120	80 - 120
1,1,1-Trichloroethane	80 - 120	80 - 120
1,1-Dichloropropene	80 - 120	80 - 120
Carbon Tetrachloride	80 - 120	80 - 123
1,2-Dichloroethane	80 - 120	80 - 120
Benzene	80 - 120	80 - 120
Trichloroethene	80 - 120	80 - 120
1,2-Dichloropropane	80 - 120	80 - 120
Bromochloromethane	80 - 120	80 - 120
Dibromomethane	80 - 120	80 - 120
2-Chloroethylvinylether	80 - 120	75 - 120
4-Methyl-2-Pentanone	80 - 120	78 - 120
cis-1,3-Dichloropropene	80 - 120	80 - 120
Toluene	80 - 120	80 - 120
trans-1,3-Dichloropropene	80 - 120	80 - 120



2-Hexanone	80 - 120	75 - 120
1,1,2-Trichloroethane	80 - 120	80 - 120
1,3-Dichloropropane	80 - 120	80 - 120
Tetrachloroethene	80 - 120	80 - 120
Dibromochloromethane	80 - 120	80 - 120
Ethylene Dibromide	80 - 120	80 - 120
Chlorobenzene	80 - 120	80 - 120
Ethylbenzene	80 - 120	80 - 121
1,1,2,2-Tetrachloroethane	80 - 120	78 - 120
m,p-Xylene	80 - 120	80 - 120
o-Xylene	80 - 120	80 - 120
Styrene	80 - 120	80 - 122
Trichlorofluoromethane	80 - 120	78 - 123
Isopropylbenzene	80 - 120	79 - 121
Bromoform	80 - 120	79 - 120
1,1,1,2-Tetrachloroethane	80 - 120	80 - 120
1,2,3-Trichloropropane	80 - 120	77 - 120
trans-1,4-Dichloro-2-butene	74 - 122	66 - 130
n-Propylbenzene	80 - 120	80 - 120
Bromobenzene	80 - 120	78 - 120
1,3,5-Trimethylbenzene	80 - 120	80 - 120
2-Chlorotoluene	80 - 120	80 - 120
4-Chlorotoluene	80 - 120	80 - 120
tert-Butylbenzene	80 - 120	80 - 121
1,2,4-Trimethylbenzene	80 - 120	80 - 120
sec-Butylbenzene	80 - 120	80 - 121
4-Isopropyltoluene	80 - 120	80 - 123
1,3-Dichlorobenzene	80 - 120	80 - 120
1,4-Dichlorobenzene	80 - 120	80 - 120
n-Butylbenzene	80 - 120	80 - 122
1,2-Dichlorobenzene	80 - 120	80 - 120
1,2-Dibromo-3-chloropropane	76 - 120	71 - 120
1,2,4-Trichlorobenzene	77 - 120	71 - 120
Hexachloro-1,3-butadiene	77 - 120	70 - 127
Naphthalene	76 - 120	70 - 120
1,2,3-Trichlorobenzene	79 - 120	74 - 120
MB/LCS Surrogate Recovery		
Dibromofluoromethane	80 - 120	(3)
d4-1,2-Dichloroethane	80 - 120	(3)
d8-Toluene	80 - 120	(3)
4-Bromofluorobenzene	80 - 120	(3)
d4-1,2-Dichlorobenzene	80 - 120	(3)
Sample Surrogate Recovery		
Dibromofluoromethane	80 - 120	(3)



d4-1,2-Dichloroethane	80 - 120	(3)
d8-Toluene	80 - 120	(3)
4-Bromofluorobenzene	80 - 120	(3)
D4-1,2-Dichlorobenzene	80 - 120	(3)

(1) Control Limits calculated using all data generated 7/1/09 through 6/30/10.

(2) ME = A marginal exceedance defined in the NELAC Standard⁽⁴⁾ as beyond the LCS-CL but still within the ME limits. ME limits are between 3 and 4 standard deviations around the mean. A maximum of four marginal exceedances are acceptable. Five or more marginal exceedances require corrective action.

(3) Marginal Exceedances not allowed for surrogate standards. A corrective action is required for each surrogate recovery outside of the control limit range.

(4) 2003 NELAC Standard (EPA/600/R-04/003), July 2003, Chapter 5, pages 251-252.

(5) Laboratory Control Sample (LCS) spike recovery control limits also used as advisory control limits for sample matrix spike (MS) analyzes. MS recovery values are advisory and not used to assess the acceptability of an analytical batch.

(6) Highlighted control limits (**bold font**) are adjusted from the calculated values as follows:

a) ARI does not use control limits < 10 for the lower limit or < 100 for the upper limit.

b) Control limits for analyzes with no separate preparation procedure are adjusted to reflect the minimum uncertainty in the calibration of the instrument allowed by the referenced analytical method.



Summary of Laboratory Control Limits Metals Analyses (All Methods & Sample Matrices)

Effective 5/1/09

Control limits are updated periodically. Assure that you have ARI's current control limits by downloading the files at the time of use. <http://www.arilabs.com/portal/downloads/ARI-CLs.zip>

Element	Matrix Spike Recovery	LCS Recovery	Replicate RPD
Aluminum	75 - 125	80 - 120	≤ 20%
Antimony	75 - 125	80 - 120	≤ 20%
Arsenic	75 - 125	80 - 120	≤ 20%
Barium	75 - 125	80 - 120	≤ 20%
Beryllium	75 - 125	80 - 120	≤ 20%
Boron	75 - 125	80 - 120	≤ 20%
Cadmium	75 - 125	80 - 120	≤ 20%
Calcium	75 - 125	80 - 120	≤ 20%
Chromium	75 - 125	80 - 120	≤ 20%
Cobalt	75 - 125	80 - 120	≤ 20%
Copper	75 - 125	80 - 120	≤ 20%
Iron	75 - 125	80 - 120	≤ 20%
Lead	75 - 125	80 - 120	≤ 20%
Magnesium	75 - 125	80 - 120	≤ 20%
Manganese	75 - 125	80 - 120	≤ 20%
Mercury	75 - 125	80 - 120	≤ 20%
Nickel	75 - 125	80 - 120	≤ 20%
Potassium	75 - 125	80 - 120	≤ 20%
Selenium	75 - 125	80 - 120	≤ 20%
Silica	75 - 125	80 - 120	≤ 20%
Silver	75 - 125	80 - 120	≤ 20%
Sodium	75 - 125	80 - 120	≤ 20%
Strontium	75 - 125	80 - 120	≤ 20%
Thallium	75 - 125	80 - 120	≤ 20%
Vanadium	75 - 125	80 - 120	≤ 20%
Zinc	75 - 125	80 - 120	≤ 20%



Spike Recovery Control Limits for Conventional Wet Chemistry

Effective 5/1/09

Control limits are updated periodically. Assure that you have ARI's current control limits by downloading the files at the time of use. <http://www.arilabs.com/portal/downloads/ARI-CLs.zip>

Sample Matrix:	ARI's Control Limits	
	Water	Soil / Sediment
Matrix Spike Recoveries	% Recovery	% Recovery
Ammonia	75 - 125	75 - 125
Bromide	75 - 125	75 - 125
Chloride	75 - 125	75 - 125
Cyanide	75 - 125	75 - 125
Ferrous Iron	75 - 125	75 - 125
Fluoride	75 - 125	75 - 125
Formaldehyde	75 - 125	75 - 125
Hexane Extractable Material	-- -- --	78 - 114
Hexavalent Chromium	75 - 125	75 - 125
Nitrate/Nitrite	75 - 125	75 - 125
Oil and Grease	75 - 125	75 - 125
Phenol	75 - 125	75 - 125
Phosphorous	75 - 125	75 - 125
Sulfate	75 - 125	75 - 125
Sulfide	75 - 125	75 - 125
Total Kjeldahl Nitrogen	75 - 125	75 - 125
Total Organic Carbon	75 - 125	75 - 125
Duplicate RPDs		
Acidity	±20%	±20%
Alkalinity	±20%	±20%
BOD	±20%	±20%
Cation Exchange	±20%	±20%
COD	±20%	±20%
Conductivity	±20%	±20%
Salinity	±20%	±20%
Solids	±20%	±20%
Turbidity	±20%	±20%

Chain of Custody Documentation

ARI Job ID: TN19, TN21

Chain of Custody Record & Laboratory Analysis Request

ARI Assigned Number: T-N-1	Turn-around Requested: Standard	Page: 1 of 1		
Client Company: AMEC	Date: 9/20/11	Ice Present? Yes		
Client Contact: John Long / Milk Bucker	No. of Coolers: 1	Cooler Temp: 0.4		
Client Project Name: FEP 2011 Shoreline Investigation	Analysis Requested			
Client Project #: 8709	Samplers: Trevor Louwrier			
Sample ID	Date	Time	Matrix	No. Containers
FEP-092011-001	9/20/11	1115	H ₂ O	5
FEP-092011-002		1300		
FEP-092011-003		1400		
FEP-092011-004		1455		
FEP-092011-005		1530		
Trip Blank		-	H ₂ O	2
<i>Handwritten note: 9/20/11</i>				
Comments/Special Instructions Metals include Al, As, Cd, Cr, Cu, Pb, Ni, Se, Th, V, Zn				
Reinstituted by: <i>M. H. Zafar</i> (Signature)	Received by: <i>J. Long</i> (Signature)	Relinquished by: <i>Trevor Louwrier</i> (Signature)		
Printed Name: Trevor Louwrier	Printed Name: Trevor Louwrier	Printed Name: Trevor Louwrier		
Company: AMEC	Company: AMEC	Company:		
Date & Time: 9/20/11 1620	Date & Time: 9-20-11 1620	Date & Time: 9-20-11 1620		

Limits of Liability: ARI will perform all requested services in accordance with appropriate methodology following ARI Standard Operating Procedures and the ARI Quality Assurance Program. This program meets standards for the industry. The total liability of ARI, its officers, agents, employees, or successors, arising out of or in connection with the requested services, shall not exceed the invoiced amount for said services. The acceptance by the client of a proposal for services by ARI release ARI from any liability in excess thereof, notwithstanding any provision to the contrary in any contract, purchase order or co-

Sample Retention Policy: All samples submitted to ARI will be appropriately discarded no sooner than 90 days after receipt or 60 days after submission of hardcopy data, whichever is longer, unless alternate retention schedules have been established by work-order or contract.



Analytical Resources, Incorporated
Analytical Chemists and Consultants

Cooler Receipt Form

ARI Client: Amec

COC No(s): _____ NA

Assigned ARI Job No: TN 19

Preliminary Examination Phase:

- Were intact, properly signed and dated custody seals attached to the outside of to cooler? YES NO
 Were custody papers included with the cooler? YES NO
 Were custody papers properly filled out (ink, signed, etc.) YES NO
 Temperature of Cooler(s) (°C) (recommended 2.0-6.0 °C for chemistry) 9.4
 If cooler temperature is out of compliance fill out form 00070F

Cooler Accepted by: TS Date: 9-20-11 Time: 1620 Temp Gun ID#: 86941619

Complete custody forms and attach all shipping documents

Log-In Phase:

- Was a temperature blank included in the cooler? YES NO
 What kind of packing material was used? ... Bubble Wrap Wet Ice Gel Packs Baggies Foam Block Paper Other:
 Was sufficient ice used (if appropriate)? YES NO
 Were all bottles sealed in individual plastic bags? YES NO
 Did all bottles arrive in good condition (unbroken)? YES NO
 Were all bottle labels complete and legible? YES NO
 Did the number of containers listed on COC match with the number of containers received? YES NO
 Did all bottle labels and tags agree with custody papers? YES NO
 Were all bottles used correct for the requested analyses? YES NO
 Do any of the analyses (bottles) require preservation? (attach preservation sheet, excluding VOCs)... NA YES NO
 Were all VOC vials free of air bubbles? NA YES NO
 Was sufficient amount of sample sent in each bottle? YES NO
 Date VOC Trip Blank was made at ARI. 9-16-11

Was Sample Split by ARI: NA YES Date/Time: _____ Equipment: _____ Split by: _____

Samples Logged by: TS Date: 9-20-11 Time: 1720

** Notify Project Manager of discrepancies or concerns **

Sample ID on Bottle	Sample ID on COC	Sample ID on Bottle	Sample ID on COC

Additional Notes, Discrepancies, & Resolutions:

FRP - 09 2011 - 003 15m11

TS 17b11

By:

Date:

<small>Small Air Bubbles ~2mm</small>	<small>Peabubbles 2-4 mm</small>	<small>LARGE Air Bubbles > 4 mm</small>	<small>Small → "sm"</small>
			<small>Peabubbles → "pb"</small>
			<small>Large → "lg"</small>
			<small>Headspace → "hs"</small>



ARI Job No: TN19

Inquiry Number: NONE
 Analysis Requested: 09/21/11
 Contact: Long, John
 Client: AMEC Geomatrix

Logged by: TS
 Sample Set Used: Yes-481
 Validatable Package: No
 Deliverables:

PC: Kelly
 VTSR: 09/20/11

Project #: 8769
 Project: FRP 2011 Shoreline Investigation
 Sample Site:
 SDG No:
 Analytical Protocol: In-house

LOGNUM ARI ID	CLIENT ID	CN >12	WAD <12	NH3 <2	COD <2	FOG <2	MET <2	PHN <2	TKN <2	NO23 <2	TOC <2	S2		AK102 Fe2+<2	DMET <2	DOC <2	FLT FLT	PARAMETER	ADJUSTED TO	LOT NUMBER	AMOUNT ADDED	DATE/BY		
												>9	<2											
11-20545 TN19A	FRP--092011-001																							
11-20546 TN19B	FRP--092011-002																							
11-20547 TN19C	FRP--092011-003																							
11-20548 TN19D	FRP--092011-004																							
11-20549 TN19E	FRP--092011-005																							

15
9-20-11
Checked By _____ Date _____



Analytical Resources, Incorporated
Analytical Chemists and Consultants

Cooler Receipt Form

ARI Client: Amec

COC No(s): _____ NA

Assigned ARI Job No: TN 21

Preliminary Examination Phase:

Were intact, properly signed and dated custody seals attached to the outside of to cooler? YES NO

Were custody papers included with the cooler? YES NO

Were custody papers properly filled out (ink, signed, etc.) YES NO

Temperature of Cooler(s) (°C) (recommended 2.0-6.0 °C for chemistry) 0.4

If cooler temperature is out of compliance fill out form 00070F

Temp Gun ID#: 96941614

Cooler Accepted by: TS Date: 9-20-11 Time: 1620

Complete custody forms and attach all shipping documents

Log-In Phase:

Was a temperature blank included in the cooler? YES NO

What kind of packing material was used? ... Bubble Wrap Wet Ice Gel Packs Baggies Foam Block Paper Other:

NA YES NO

Was sufficient ice used (if appropriate)? YES NO

Were all bottles sealed in individual plastic bags? YES NO

Did all bottles arrive in good condition (unbroken)? YES NO

Were all bottle labels complete and legible? YES NO

Did the number of containers listed on COC match with the number of containers received? YES NO

Did all bottle labels and tags agree with custody papers? YES NO

Were all bottles used correct for the requested analyses? YES NO

Do any of the analyses (bottles) require preservation? (attach preservation sheet, excluding VOCs)... NA YES NO

Were all VOC vials free of air bubbles? NA YES NO

Was sufficient amount of sample sent in each bottle? YES NO

Date VOC Trip Blank was made at ARI: 9-16-11

Was Sample Split by ARI: NO YES Date/Time: _____ Equipment: _____ Split by: _____

Samples Logged by: TS Date: 9-20-11 Time: 1720

*** Notify Project Manager of discrepancies or concerns ***

Sample ID on Bottle	Sample ID on COC	Sample ID on Bottle	Sample ID on COC

Additional Notes, Discrepancies, & Resolutions:

FRP - 09 2011 - 003 15mll

1B 1pb11

By:

Date:

Small Air Bubbles ~2mm * * * *	Peabubbles 2-4 mm * * * *	LARGE Air Bubbles > 4 mm * * * *	Small → "sm" Peabubbles → "pb" Large → "lg" Headspace → "hs"

PRESERVATION VERIFICATION 09/20/11

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Inquiry Number: NONE
Analysis Requested: 09/21/11
Contact: Long, John
Client: AMEC Geomatix
Logged by: TS
Sample Set Used: Yes-481
Validatable Package: No
Deliverables:

ARI Job No: TN19

Project #: 8769
Project: FRP 2011 Shoreline Investigation
Sample Site:
SDG No:
Analytical Protocol: In-house

THIS IS A PUBLIC RECORD

Case Narrative, Data Qualifiers, Control Limits

ARI Job ID: TN19, TN21

Case Narrative

AMEC/Geomatrix

Client Project: Former Rhone Poulenc- 8769 Shoreline Investigation

ARI Job Numbers: TN19, TN21

October 13, 2011

Sample Receipt:

Please find enclosed the original Chain-of-Custody (COC) record, sample receipt documentation, and analytical results for the project referenced above. Analytical Resources, Inc. accepted five water samples and one trip blank in good condition on 9/20/11. Please see the enclosed Cooler Receipt Form for further details.

Volatiles by 8260C

The samples were analyzed on 9/21/11 and 9/26/11 within the method recommended holding times.

Initial calibration (s): All analytes of interest were within method acceptance criteria.

Continuing calibration (s): The continuing calibration (CCAL) on 9/21/11 fell outside the 20% control limit low for Hexachlorobutadiene. All detected results associated with this CCAL have been flagged with a "Q" qualifier.

The CCAL on 9/26/11 fell outside the 20% control limit low for Acetone, Methyl Iodide, and trans-1,4-Dichloro-2-butene. All detected results associated with this CCAL have been flagged with a "Q" qualifier.

LCS/LCSD/RPDs: The LCSD percent recovery of Isopropylbenzene was outside the control limits high for LCS-092111

The LCS percent recovery of 2-Chloroethylvinylether fell outside the control limits low for LCS-092611.

Surrogates: All surrogate recoveries were within control limits.

Method Blank (s): The method blanks were free of contamination.

Samples: Samples FRP-092011-001 and FRP-092011-005 had a pH between 6 and 8.

Metals Analysis (6010, 200.8 and 7000 series)

The samples were digested on 9/22/11 - within the method recommended holding time and analyzed between 9/27/11 and 10/7/11.

Initial calibration (s): All analytes of interest were within method acceptance criteria.

Continuing calibration (s): All analytes of interest were within method acceptance criteria.

Case Narrative

AMEC/Geomatrix

Client Project: Former Rhone Poulenc- 8769 Shoreline Investigation

ARI Job Numbers: TN19, TN21

October 13, 2011

LCS (s): The percent recoveries were within control limits.

Method Blank (s): The method blanks were free of contamination.

Samples: There were no anomalies associated with these samples.

Matrix spike/ Sample Duplicate RPD(s): The matrix spike percent recovery of selenium fell outside the control limits high for sample **FRP-092011-001**. A post digestion spike was performed and the recovery was within control limits. All relevant data have been flagged with an "N" qualifier on the appropriate Form V.

The matrix spike percent recovery of mercury fell outside the control limits low for sample **FRP-092011-001**. All relevant data have been flagged with an "N" qualifier on the appropriate Form V.

The duplicate RPDs were within control limits.

pH by method 150.1:

The samples were analyzed on 9/20/11 within method recommended holding time.

LCS (s): The LCS percent recovery was within control limits.

Sample Replicate RPD (s): The RPD is in control.

Sample ID Cross Reference Report

ARI Job No: TN19
Client: AMEC Geomatrix
Project Event: 8769
Project Name: FRP 2011 Shoreline Investigation

Sample ID	ARI Lab ID	ARI LIMS ID	Matrix	Sample Date/Time	VTSR
1. FRP-092011-001	TN19A	11-20545	Water	09/20/11 11:15	09/20/11 16:20
2. FRP-092011-002	TN19B	11-20546	Water	09/20/11 13:00	09/20/11 16:20
3. FRP-092011-003	TN19C	11-20547	Water	09/20/11 14:00	09/20/11 16:20
4. FRP-092011-004	TN19D	11-20548	Water	09/20/11 14:55	09/20/11 16:20
5. FRP-092011-005	TN19E	11-20549	Water	09/20/11 15:30	09/20/11 16:20
6. Trip Blanks	TN19F	11-20550	Water	09/20/11	09/20/11 16:20

Printed 09/20/11

Sample ID Cross Reference Report

ARI Job No: TN21
Client: AMEC Geomatrix
Project Event: 8769
Project Name: FRP 2011 Shoreline Investigation

Sample ID	ARI Lab ID	ARI LIMS ID	Matrix	Sample Date/Time	VTSR
1. FRP-092011-001	TN21A	11-20551	Water	09/20/11 11:15	09/20/11 16:20
2. FRP-092011-002	TN21B	11-20552	Water	09/20/11 13:00	09/20/11 16:20
3. FRP-092011-003	TN21C	11-20553	Water	09/20/11 14:00	09/20/11 16:20
4. FRP-092011-004	TN21D	11-20554	Water	09/20/11 14:55	09/20/11 16:20
5. FRP-092011-005	TN21E	11-20555	Water	09/20/11 15:30	09/20/11 16:20

Printed 09/20/11

TN19:00012



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Data Reporting Qualifiers

Effective 2/14/2011

Inorganic Data

- U Indicates that the target analyte was not detected at the reported concentration
- * Duplicate RPD is not within established control limits
- B Reported value is less than the CRDL but \geq the Reporting Limit
- N Matrix Spike recovery not within established control limits
- NA Not Applicable, analyte not spiked
- H The natural concentration of the spiked element is so much greater than the concentration spiked that an accurate determination of spike recovery is not possible
- L Analyte concentration is \leq 5 times the Reporting Limit and the replicate control limit defaults to ± 1 RL instead of the normal 20% RPD

Organic Data

- U Indicates that the target analyte was not detected at the reported concentration
- * Flagged value is not within established control limits
- B Analyte detected in an associated Method Blank at a concentration greater than one-half of ARI's Reporting Limit or 5% of the regulatory limit or 5% of the analyte concentration in the sample.
- J Estimated concentration when the value is less than ARI's established reporting limits
- D The spiked compound was not detected due to sample extract dilution
- E Estimated concentration calculated for an analyte response above the valid instrument calibration range. A dilution is required to obtain an accurate quantification of the analyte.
- Q Indicates a detected analyte with an initial or continuing calibration that does not meet established acceptance criteria (<20%RSD, <20%Drift or minimum RRF).



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- S Indicates an analyte response that has saturated the detector. The calculated concentration is not valid; a dilution is required to obtain valid quantification of the analyte
- NA The flagged analyte was not analyzed for
- NR Spiked compound recovery is not reported due to chromatographic interference
- NS The flagged analyte was not spiked into the sample
- M Estimated value for an analyte detected and confirmed by an analyst but with low spectral match parameters. This flag is used only for GC-MS analyses
- M2 The sample contains PCB congeners that do not match any standard Aroclor pattern. The PCBs are identified and quantified as the Aroclor whose pattern most closely matches that of the sample. The reported value is an estimate.
- N The analysis indicates the presence of an analyte for which there is presumptive evidence to make a "tentative identification"
- Y The analyte is not detected at or above the reported concentration. The reporting limit is raised due to chromatographic interference. The Y flag is equivalent to the U flag with a raised reporting limit.
- EMPC Estimated Maximum Possible Concentration (EMPC) defined in EPA Statement of Work DLM02.2 as a value "calculated for 2,3,7,8-substituted isomers for which the quantitation and /or confirmation ion(s) has signal to noise in excess of 2.5, but does not meet identification criteria" **(Dioxin/Furan analysis only)**
- C The analyte was positively identified on only one of two chromatographic columns. Chromatographic interference prevented a positive identification on the second column
- P The analyte was detected on both chromatographic columns but the quantified values differ by ≥40% RPD with no obvious chromatographic interference
- X Analyte signal includes interference from polychlorinated diphenyl ethers. **(Dioxin/Furan analysis only)**
- Z Analyte signal includes interference from the sample matrix or perfluorokerosene ions. **(Dioxin/Furan analysis only)**



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Geotechnical Data

- A The total of all fines fractions. This flag is used to report total fines when only sieve analysis is requested and balances total grain size with sample weight.
- F Samples were frozen prior to particle size determination
- SM Sample matrix was not appropriate for the requested analysis. This normally refers to samples contaminated with an organic product that interferes with the sieving process and/or moisture content, porosity and saturation calculations
- SS Sample did not contain the proportion of "fines" required to perform the pipette portion of the grain size analysis
- W Weight of sample in some pipette aliquots was below the level required for accurate weighting

SURR SOLUTIONS

LABEL SOLN ID TEST CONC. UG/ML SOLVENT EXP.

LCS SOLUTIONS

LABL	SOLN ID	TEST	CONC. UG/ML	SOLVENT	EXP.
1	1888-2	PCB 1660	20	ACETONE	08/30/12
2#	NA	BCOC PEST	10	ACETONE	NA
3	1885-1	PEST	01/02/10	ACETONE	12/15/11
4	1885-2	LOW PEST	.1/.2/1	ACETONE	12/15/11
5	1779-1	EPH	1500	MECL2	11/11/11
6	1791-5	PCP	12.5/125	ACETONE	12/10/11
7	1888-1	ABN	100	MEOH	08/30/12
8	1785-3	TBT	2.5	MECL2	11/27/11
9	1786-3	PORE TBT	.125/.25	MECL2	11/27/11
10					
11	1860-4	TPHD	15000	ACETONE	05/12/12
12					
13	1838-4	LOW PCB	2	ACETONE	01/31/12
14					
15	1814-2	SIM PNA	15/75	MEOH	01/04/12
16	1879-3	1,4-DIOXANE	100	MEOH	02/05/12
17	1869-4	1248 PCB	10	ACETONE	06/14/12
18	1814-3	LOW SIM PNA	1.5	ACETONE	01/04/12
19	1873-2	AK103	7500	ACETONE	01/02/12
20	1886-4	PNA	100	ACETONE	01/07/12
21	1874-3	SKY/BHT	100	MEOH	01/14/12
22	1864-3	HERB	02 to 2500	MEOH	12/03/11
23	1887-2	EXTRA PNA	15	ACETONE	08/25/12
24					
25#	NA	DIPHENYL	100	MEOH	NA
26	1869-1	OP-PEST	25	MEOH	10/01/11
27	NA	STEROLS	200	MEOH	NA
28#	1807-1	ADD. PEST	2	ACETONE	08/31/11
29#	NA	DECANES	100	MEOH	NA

LCS SOLUTIONS

30	NA	EDB/DBCP	0.2	MEOH	NA
31	1835-2	TERPINEOL	100	MEOH	09/02/11
32	1876-1	GUAIACOL	50-200	ACETONE	01/05/12
33	NA	RETENE	100	MEOH	NA
34	1867-3	CONGENERS	0.5	ACETONE	03/14/12
35	1875-3	ALKYL PNA A	10	MEOH	07/18/12
36	NA	ALKYL PNA B	10	MEOH	NA
37	1773-1	CAR/PERY	100	ACETONE	10/14/11
38	1872-2	ABN ACID	200-450	MEOH	12/29/11
39	1853-4	BENZIDINE	500	MEOH	04/30/12
40	1851-3	PBDE	0.5	MEOH	04/22/12
50	1900-1	FULL RESIN	250	ACETONE	08/12/12
51	1772-1	DDTS	0.01	ACETONE	04/24/11
52	NA	1232 PCB	20	ACETONE	NA
53	1852-2	DALAPON	50	MEOH	12/03/11
54	1753-1	T-CHLORDANE	10	ACETONE	07/21/11
55	1753-2	TOXAPHENE	50	ACETONE	07/21/11
56	1874-1	ABN BASE	50-200	MEOH	01/05/12
#=PROJECT SPECIFIC SOLUTION					
*=REVERIFIED SOLUTION					

Volatile Analysis
Report and Summary QC Forms

ARI Job ID: TN19, TN21

ORGANICS ANALYSIS DATA SHEET

Volatile by Purge & Trap GC/MS-Method SW8260C
Page 1 of 2
**ANALYTICAL
RESOURCES
INCORPORATED**


Sample ID: FRP-092011-001

SAMPLE

Lab Sample ID: TN19A
 LIMS ID: 11-20545
 Matrix: Water
 Data Release Authorized: *MW*
 Reported: 10/03/11

QC Report No: TN19-AMEC Geomatrix
 Project: FRP 2011 Shoreline Investigation
 8769
 Date Sampled: 09/20/11
 Date Received: 09/20/11

Instrument/Analyst: NT5/PAB
 Date Analyzed: 09/21/11 17:22

Sample Amount: 2.00 mL
 Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.49	2.5	< 2.5 U
74-83-9	Bromomethane	0.22	5.0	< 5.0 U
75-01-4	Vinyl Chloride	0.38	1.0	< 1.0 U
75-00-3	Chloroethane	0.76	1.0	< 1.0 U
75-09-2	Methylene Chloride	2.0	2.5	< 2.5 U
67-64-1	Acetone	3.6	25	< 25 U
75-15-0	Carbon Disulfide	0.44	1.0	< 1.0 U
75-35-4	1,1-Dichloroethene	0.46	1.0	< 1.0 U
75-34-3	1,1-Dichloroethane	0.26	1.0	< 1.0 U
156-60-5	trans-1,2-Dichloroethene	0.42	1.0	< 1.0 U
156-59-2	cis-1,2-Dichloroethene	0.50	1.0	< 1.0 U
67-66-3	Chloroform	0.40	1.0	< 1.0 U
107-06-2	1,2-Dichloroethane	0.38	1.0	< 1.0 U
78-93-3	2-Butanone	4.0	25	< 25 U
71-55-6	1,1,1-Trichloroethane	0.44	1.0	< 1.0 U
56-23-5	Carbon Tetrachloride	0.38	1.0	< 1.0 U
108-05-4	Vinyl Acetate	0.34	5.0	< 5.0 U
75-27-4	Bromodichloromethane	0.26	1.0	< 1.0 U
78-87-5	1,2-Dichloropropane	0.46	1.0	< 1.0 U
10061-01-5	cis-1,3-Dichloropropene	0.29	1.0	< 1.0 U
79-01-6	Trichloroethene	0.38	1.0	< 1.0 U
124-48-1	Dibromochloromethane	0.45	1.0	< 1.0 U
79-00-5	1,1,2-Trichloroethane	0.18	1.0	< 1.0 U
71-43-2	Benzene	0.28	1.0	< 1.0 U
10061-02-6	trans-1,3-Dichloropropene	0.30	1.0	< 1.0 U
110-75-8	2-Chloroethylvinylether	0.43	5.0	< 5.0 U
75-25-2	Bromoform	0.35	1.0	< 1.0 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	1.9	25	< 25 U
591-78-6	2-Hexanone	1.6	25	< 25 U
127-18-4	Tetrachloroethene	0.44	1.0	< 1.0 U
79-34-5	1,1,2,2-Tetrachloroethane	0.34	1.0	< 1.0 U
108-88-3	Toluene	0.28	1.0	< 1.0 U
108-90-7	Chlorobenzene	0.21	1.0	< 1.0 U
100-41-4	Ethylbenzene	0.47	1.0	< 1.0 U
100-42-5	Styrene	0.33	1.0	< 1.0 U
75-69-4	Trichlorofluoromethane	0.46	1.0	< 1.0 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.54	1.0	< 1.0 U
179601-23-1	m,p-Xylene	0.72	2.0	< 2.0 U
95-47-6	o-Xylene	0.28	1.0	< 1.0 U
95-50-1	1,2-Dichlorobenzene	0.28	1.0	< 1.0 U
541-73-1	1,3-Dichlorobenzene	0.20	1.0	< 1.0 U
106-46-7	1,4-Dichlorobenzene	0.28	1.0	< 1.0 U
107-02-8	Acrolein	1.5	25	< 25 U
74-88-4	Methyl Iodide	0.20	5.0	< 5.0 U
74-96-4	Bromoethane	0.45	1.0	< 1.0 U
107-13-1	Acrylonitrile	0.92	5.0	< 5.0 U
563-58-6	1,1-Dichloropropene	0.46	1.0	< 1.0 U
74-95-3	Dibromomethane	0.40	1.0	< 1.0 U
630-20-6	1,1,1,2-Tetrachloroethane	0.34	1.0	< 1.0 U
96-12-8	1,2-Dibromo-3-chloropropane	1.1	2.5	< 2.5 U

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2

Sample ID: FRP-092011-001
SAMPLE

Lab Sample ID: TN19A
LIMS ID: 11-20545
Matrix: Water
Date Analyzed: 09/21/11 17:22

QC Report No: TN19-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	1.1	2.5	< 2.5 U
110-57-6	trans-1,4-Dichloro-2-butene	1.2	5.0	< 5.0 U
108-67-8	1,3,5-Trimethylbenzene	0.32	1.0	< 1.0 U
95-63-6	1,2,4-Trimethylbenzene	0.29	1.0	< 1.0 U
87-68-3	Hexachlorobutadiene	0.56	2.5	< 2.5 U
106-93-4	Ethylene Dibromide	0.38	1.0	< 1.0 U
74-97-5	Bromochloromethane	0.34	1.0	< 1.0 U
594-20-7	2,2-Dichloropropane	0.42	1.0	< 1.0 U
142-28-9	1,3-Dichloropropane	0.10	1.0	< 1.0 U
98-82-8	Isopropylbenzene	0.31	1.0	< 1.0 U
103-65-1	n-Propylbenzene	0.40	1.0	< 1.0 U
108-86-1	Bromobenzene	0.26	1.0	< 1.0 U
95-49-8	2-Chlorotoluene	0.21	1.0	< 1.0 U
106-43-4	4-Chlorotoluene	0.36	1.0	< 1.0 U
98-06-6	tert-Butylbenzene	0.30	1.0	< 1.0 U
135-98-8	sec-Butylbenzene	0.38	1.0	< 1.0 U
99-87-6	4-Isopropyltoluene	0.38	1.0	< 1.0 U
104-51-8	n-Butylbenzene	0.54	1.0	< 1.0 U
120-82-1	1,2,4-Trichlorobenzene	0.50	2.5	< 2.5 U
91-20-3	Naphthalene	0.35	2.5	< 2.5 U
87-61-6	1,2,3-Trichlorobenzene	0.44	2.5	< 2.5 U

Reported in $\mu\text{g/L}$ (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	102%
d8-Toluene	100%
Bromofluorobenzene	93.8%
d4-1,2-Dichlorobenzene	98.1%

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

ORGANICS ANALYSIS DATA SHEET

Volatile by Purge & Trap GC/MS-Method SW8260C
Page 1 of 2

Sample ID: FRP-092011-002

SAMPLE

Lab Sample ID: TN19B

LIMS ID: 11-20546

Matrix: Water

Data Release Authorized: *TWW*

Reported: 10/03/11

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: 09/20/11

Date Received: 09/20/11

Instrument/Analyst: NT2/PAB

Date Analyzed: 09/26/11 14:42

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.10	0.5	< 0.5 U
74-83-9	Bromomethane	0.04	1.0	< 1.0 U
75-01-4	Vinyl Chloride	0.08	0.2	< 0.2 U
75-00-3	Chloroethane	0.15	0.2	< 0.2 U
75-09-2	Methylene Chloride	0.39	0.5	< 0.5 U
67-64-1	Acetone	0.72	5.0	< 5.0 U
75-15-0	Carbon Disulfide	0.09	0.2	< 0.2 U
75-35-4	1,1-Dichloroethene	0.09	0.2	< 0.2 U
75-34-3	1,1-Dichloroethane	0.05	0.2	< 0.2 U
156-60-5	trans-1,2-Dichloroethene	0.08	0.2	< 0.2 U
156-59-2	cis-1,2-Dichloroethene	0.10	0.2	< 0.2 U
67-66-3	Chloroform	0.08	0.2	< 0.2 U
107-06-2	1,2-Dichloroethane	0.08	0.2	< 0.2 U
78-93-3	2-Butanone	0.81	5.0	< 5.0 U
71-55-6	1,1,1-Trichloroethane	0.09	0.2	< 0.2 U
56-23-5	Carbon Tetrachloride	0.08	0.2	< 0.2 U
108-05-4	Vinyl Acetate	0.07	1.0	< 1.0 U
75-27-4	Bromodichloromethane	0.05	0.2	< 0.2 U
78-87-5	1,2-Dichloropropane	0.09	0.2	< 0.2 U
10061-01-5	cis-1,3-Dichloropropene	0.06	0.2	< 0.2 U
79-01-6	Trichloroethene	0.08	0.2	< 0.2 U
124-48-1	Dibromochloromethane	0.09	0.2	< 0.2 U
79-00-5	1,1,2-Trichloroethane	0.04	0.2	< 0.2 U
71-43-2	Benzene	0.06	0.2	< 0.2 U
10061-02-6	trans-1,3-Dichloropropene	0.06	0.2	< 0.2 U
110-75-8	2-Chloroethylvinylether	0.09	1.0	< 1.0 U
75-25-2	Bromoform	0.07	0.2	< 0.2 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	0.38	5.0	< 5.0 U
591-78-6	2-Hexanone	0.31	5.0	< 5.0 U
127-18-4	Tetrachloroethene	0.09	0.2	< 0.2 U
79-34-5	1,1,2,2-Tetrachloroethane	0.07	0.2	< 0.2 U
108-88-3	Toluene	0.06	0.2	< 0.2 U
108-90-7	Chlorobenzene	0.04	0.2	< 0.2 U
100-41-4	Ethylbenzene	0.09	0.2	< 0.2 U
100-42-5	Styrene	0.07	0.2	< 0.2 U
75-69-4	Trichlorofluoromethane	0.09	0.2	< 0.2 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	0.11	0.2	< 0.2 U
179601-23-1	m,p-Xylene	0.14	0.4	< 0.4 U
95-47-6	o-Xylene	0.06	0.2	< 0.2 U
95-50-1	1,2-Dichlorobenzene	0.06	0.2	< 0.2 U
541-73-1	1,3-Dichlorobenzene	0.04	0.2	< 0.2 U
106-46-7	1,4-Dichlorobenzene	0.06	0.2	< 0.2 U
107-02-8	Acrolein	0.29	5.0	< 5.0 U
74-88-4	Methyl Iodide	0.04	1.0	< 1.0 U
74-96-4	Bromoethane	0.09	0.2	< 0.2 U
107-13-1	Acrylonitrile	0.18	1.0	< 1.0 U
563-58-6	1,1-Dichloropropene	0.09	0.2	< 0.2 U
74-95-3	Dibromomethane	0.08	0.2	< 0.2 U
630-20-6	1,1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
96-12-8	1,2-Dibromo-3-chloropropane	0.21	0.5	< 0.5 U

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C

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**ANALYTICAL
RESOURCES
INCORPORATED**


Sample ID: FRP-092011-002

SAMPLE

Lab Sample ID: TN19B

LIMS ID: 11-20546

Matrix: Water

Date Analyzed: 09/26/11 14:42

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	0.23	0.5	< 0.5 U
110-57-6	trans-1,4-Dichloro-2-butene	0.24	1.0	< 1.0 U
108-67-8	1,3,5-Trimethylbenzene	0.06	0.2	< 0.2 U
95-63-6	1,2,4-Trimethylbenzene	0.06	0.2	< 0.2 U
87-68-3	Hexachlorobutadiene	0.11	0.5	< 0.5 U
106-93-4	Ethylene Dibromide	0.08	0.2	< 0.2 U
74-97-5	Bromo-chloromethane	0.07	0.2	< 0.2 U
594-20-7	2,2-Dichloropropane	0.08	0.2	< 0.2 U
142-28-9	1,3-Dichloropropane	0.02	0.2	< 0.2 U
98-82-8	Isopropylbenzene	0.06	0.2	< 0.2 U
103-65-1	n-Propylbenzene	0.08	0.2	< 0.2 U
108-86-1	Bromobenzene	0.05	0.2	< 0.2 U
95-49-8	2-Chlorotoluene	0.04	0.2	< 0.2 U
106-43-4	4-Chlorotoluene	0.07	0.2	< 0.2 U
98-06-6	tert-Butylbenzene	0.06	0.2	< 0.2 U
135-98-8	sec-Butylbenzene	0.08	0.2	< 0.2 U
99-87-6	4-Isopropyltoluene	0.08	0.2	< 0.2 U
104-51-8	n-Butylbenzene	0.11	0.2	< 0.2 U
120-82-1	1,2,4-Trichlorobenzene	0.10	0.5	< 0.5 U
91-20-3	Naphthalene	0.07	0.5	< 0.5 U
87-61-6	1,2,3-Trichlorobenzene	0.09	0.5	< 0.5 U

Reported in µg/L (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	93.9%
d8-Toluene	100%
Bromo-fluorobenzene	96.3%
d4-1,2-Dichlorobenzene	102%

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

ORGANICS ANALYSIS DATA SHEET

Volatile by Purge & Trap GC/MS-Method SW8260C

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**ANALYTICAL
RESOURCES
INCORPORATED**


Sample ID: FRP-092011-003

SAMPLE

Lab Sample ID: TN19C

LIMS ID: 11-20547

Matrix: Water

Data Release Authorized: **YWW**

Reported: 10/03/11

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: 09/20/11

Date Received: 09/20/11

Instrument/Analyst: NT2/PAB

Date Analyzed: 09/26/11 15:09

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.10	0.5	< 0.5 U
74-83-9	Bromomethane	0.04	1.0	< 1.0 U
75-01-4	Vinyl Chloride	0.08	0.2	< 0.2 U
75-00-3	Chloroethane	0.15	0.2	< 0.2 U
75-09-2	Methylene Chloride	0.39	0.5	< 0.5 U
67-64-1	Acetone	0.72	5.0	< 5.0 U
75-15-0	Carbon Disulfide	0.09	0.2	< 0.2 U
75-35-4	1,1-Dichloroethene	0.09	0.2	< 0.2 U
75-34-3	1,1-Dichloroethane	0.05	0.2	< 0.2 U
156-60-5	trans-1,2-Dichloroethene	0.08	0.2	< 0.2 U
156-59-2	cis-1,2-Dichloroethene	0.10	0.2	< 0.2 U
67-66-3	Chloroform	0.08	0.2	< 0.2 U
107-06-2	1,2-Dichloroethane	0.08	0.2	< 0.2 U
78-93-3	2-Butanone	0.81	5.0	< 5.0 U
71-55-6	1,1,1-Trichloroethane	0.09	0.2	< 0.2 U
56-23-5	Carbon Tetrachloride	0.08	0.2	< 0.2 U
108-05-4	Vinyl Acetate	0.07	1.0	< 1.0 U
75-27-4	Bromodichloromethane	0.05	0.2	< 0.2 U
78-87-5	1,2-Dichloropropane	0.09	0.2	< 0.2 U
10061-01-5	cis-1,3-Dichloropropene	0.06	0.2	< 0.2 U
79-01-6	Trichloroethene	0.08	0.2	< 0.2 U
124-48-1	Dibromochloromethane	0.09	0.2	< 0.2 U
79-00-5	1,1,2-Trichloroethane	0.04	0.2	< 0.2 U
71-43-2	Benzene	0.06	0.2	< 0.2 U
10061-02-6	trans-1,3-Dichloropropene	0.06	0.2	< 0.2 U
110-75-8	2-Chloroethylvinylether	0.09	1.0	< 1.0 U
75-25-2	Bromoform	0.07	0.2	< 0.2 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	0.38	5.0	< 5.0 U
591-78-6	2-Hexanone	0.31	5.0	< 5.0 U
127-18-4	Tetrachloroethene	0.09	0.2	< 0.2 U
79-34-5	1,1,2,2-Tetrachloroethane	0.07	0.2	< 0.2 U
108-88-3	Toluene	0.06	0.2	< 0.2 U
108-90-7	Chlorobenzene	0.04	0.2	< 0.2 U
100-41-4	Ethylbenzene	0.09	0.2	< 0.2 U
100-42-5	Styrene	0.07	0.2	< 0.2 U
75-69-4	Trichlorofluoromethane	0.09	0.2	< 0.2 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoro	0.11	0.2	< 0.2 U
179601-23-1	m,p-Xylene	0.14	0.4	< 0.4 U
95-47-6	o-Xylene	0.06	0.2	< 0.2 U
95-50-1	1,2-Dichlorobenzene	0.06	0.2	< 0.2 U
541-73-1	1,3-Dichlorobenzene	0.04	0.2	< 0.2 U
106-46-7	1,4-Dichlorobenzene	0.06	0.2	< 0.2 U
107-02-8	Acrolein	0.29	5.0	< 5.0 U
74-88-4	Methyl Iodide	0.04	1.0	< 1.0 U
74-96-4	Bromoethane	0.09	0.2	< 0.2 U
107-13-1	Acrylonitrile	0.18	1.0	< 1.0 U
563-58-6	1,1-Dichloropropene	0.09	0.2	< 0.2 U
74-95-3	Dibromomethane	0.08	0.2	< 0.2 U
630-20-6	1,1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
96-12-8	1,2-Dibromo-3-chloropropane	0.21	0.5	< 0.5 U

OF
2011

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
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Sample ID: FRP-092011-003
SAMPLE

Lab Sample ID: TN19C
LIMS ID: 11-20547
Matrix: Water
Date Analyzed: 09/26/11 15:09

QC Report No: TN19-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	0.23	0.5	< 0.5 U
110-57-6	trans-1,4-Dichloro-2-butene	0.24	1.0	< 1.0 U
108-67-8	1,3,5-Trimethylbenzene	0.06	0.2	< 0.2 U
95-63-6	1,2,4-Trimethylbenzene	0.06	0.2	< 0.2 U
87-68-3	Hexachlorobutadiene	0.11	0.5	< 0.5 U
106-93-4	Ethylene Dibromide	0.08	0.2	< 0.2 U
74-97-5	Bromo-chloromethane	0.07	0.2	< 0.2 U
594-20-7	2,2-Dichloropropane	0.08	0.2	< 0.2 U
142-28-9	1,3-Dichloropropane	0.02	0.2	< 0.2 U
98-82-8	Isopropylbenzene	0.06	0.2	< 0.2 U
103-65-1	n-Propylbenzene	0.08	0.2	< 0.2 U
108-86-1	Bromobenzene	0.05	0.2	< 0.2 U
95-49-8	2-Chlorotoluene	0.04	0.2	< 0.2 U
106-43-4	4-Chlorotoluene	0.07	0.2	< 0.2 U
98-06-6	tert-Butylbenzene	0.06	0.2	< 0.2 U
135-98-8	sec-Butylbenzene	0.08	0.2	< 0.2 U
99-87-6	4-Isopropyltoluene	0.08	0.2	< 0.2 U
104-51-8	n-Butylbenzene	0.11	0.2	< 0.2 U
120-82-1	1,2,4-Trichlorobenzene	0.10	0.5	< 0.5 U
91-20-3	Naphthalene	0.07	0.5	< 0.5 U
87-61-6	1,2,3-Trichlorobenzene	0.09	0.5	< 0.5 U

Reported in $\mu\text{g/L}$ (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	93.9%
d8-Toluene	99.5%
Bromo-fluorobenzene	95.7%
d4-1,2-Dichlorobenzene	102%

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

ORGANICS ANALYSIS DATA SHEET

Volatile by Purge & Trap GC/MS-Method SW8260C

Page 1 of 2



Sample ID: FRP-092011-004

SAMPLE

Lab Sample ID: TN19D

LIMS ID: 11-20548

Matrix: Water

Data Release Authorized: *WW*

Reported: 10/03/11

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: 09/20/11

Date Received: 09/20/11

Instrument/Analyst: NT2/PAB

Date Analyzed: 09/26/11 15:36

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.10	0.5	< 0.5 U
74-83-9	Bromomethane	0.04	1.0	< 1.0 U
75-01-4	Vinyl Chloride	0.08	0.2	< 0.2 U
75-00-3	Chloroethane	0.15	0.2	< 0.2 U
75-09-2	Methylene Chloride	0.39	0.5	< 0.5 U
67-64-1	Acetone	0.72	5.0	< 5.0 U <i>J</i>
75-15-0	Carbon Disulfide	0.09	0.2	< 0.2 U
75-35-4	1,1-Dichloroethene	0.09	0.2	< 0.2 U
75-34-3	1,1-Dichloroethane	0.05	0.2	< 0.2 U
156-60-5	trans-1,2-Dichloroethene	0.08	0.2	< 0.2 U
156-59-2	cis-1,2-Dichloroethene	0.10	0.2	< 0.2 U
67-66-3	Chloroform	0.08	0.2	< 0.2 U
107-06-2	1,2-Dichloroethane	0.08	0.2	< 0.2 U
78-93-3	2-Butanone	0.81	5.0	< 5.0 U
71-55-6	1,1,1-Trichloroethane	0.09	0.2	< 0.2 U
56-23-5	Carbon Tetrachloride	0.08	0.2	< 0.2 U
108-05-4	Vinyl Acetate	0.07	1.0	< 1.0 U
75-27-4	Bromodichloromethane	0.05	0.2	< 0.2 U
78-87-5	1,2-Dichloropropane	0.09	0.2	< 0.2 U
10061-01-5	cis-1,3-Dichloropropene	0.06	0.2	< 0.2 U
79-01-6	Trichloroethene	0.08	0.2	< 0.2 U
124-48-1	Dibromochloromethane	0.09	0.2	< 0.2 U
79-00-5	1,1,2-Trichloroethane	0.04	0.2	< 0.2 U
71-43-2	Benzene	0.06	0.2	< 0.2 U
10061-02-6	trans-1,3-Dichloropropene	0.06	0.2	< 0.2 U
110-75-8	2-Chloroethylvinylether	0.09	1.0	< 1.0 U
75-25-2	Bromoform	0.07	0.2	< 0.2 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	0.38	5.0	< 5.0 U
591-78-6	2-Hexanone	0.31	5.0	< 5.0 U
127-18-4	Tetrachloroethene	0.09	0.2	< 0.2 U
79-34-5	1,1,2,2-Tetrachloroethane	0.07	0.2	< 0.2 U
108-88-3	Toluene	0.06	0.2	< 0.2 U
108-90-7	Chlorobenzene	0.04	0.2	< 0.2 U
100-41-4	Ethylbenzene	0.09	0.2	< 0.2 U
100-42-5	Styrene	0.07	0.2	< 0.2 U
75-69-4	Trichlorofluoromethane	0.09	0.2	< 0.2 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoro	0.11	0.2	< 0.2 U
179601-23-1	m,p-Xylene	0.14	0.4	< 0.4 U
95-47-6	o-Xylene	0.06	0.2	< 0.2 U
95-50-1	1,2-Dichlorobenzene	0.06	0.2	< 0.2 U
541-73-1	1,3-Dichlorobenzene	0.04	0.2	< 0.2 U
106-46-7	1,4-Dichlorobenzene	0.06	0.2	< 0.2 U
107-02-8	Acrolein	0.29	5.0	< 5.0 U
74-88-4	Methyl Iodide	0.04	1.0	< 1.0 U <i>J</i>
74-96-4	Bromoethane	0.09	0.2	< 0.2 U
107-13-1	Acrylonitrile	0.18	1.0	< 1.0 U
563-58-6	1,1-Dichloropropene	0.09	0.2	< 0.2 U
74-95-3	Dibromomethane	0.08	0.2	< 0.2 U
630-20-6	1,1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
96-12-8	1,2-Dibromo-3-chloropropane	0.21	0.5	< 0.5 U

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2

Sample ID: FRP-092011-004

SAMPLE

Lab Sample ID: TN19D
LIMS ID: 11-20548
Matrix: Water
Date Analyzed: 09/26/11 15:36

QC Report No: TN19-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	0.23	0.5	< 0.5 U
110-57-6	trans-1,4-Dichloro-2-butene	0.24	1.0	< 1.0 U
108-67-8	1,3,5-Trimethylbenzene	0.06	0.2	< 0.2 U
95-63-6	1,2,4-Trimethylbenzene	0.06	0.2	< 0.2 U
87-68-3	Hexachlorobutadiene	0.11	0.5	< 0.5 U
106-93-4	Ethylene Dibromide	0.08	0.2	< 0.2 U
74-97-5	Bromo-chloromethane	0.07	0.2	< 0.2 U
594-20-7	2,2-Dichloropropane	0.08	0.2	< 0.2 U
142-28-9	1,3-Dichloropropane	0.02	0.2	< 0.2 U
98-82-8	Isopropylbenzene	0.06	0.2	< 0.2 U
103-65-1	n-Propylbenzene	0.08	0.2	< 0.2 U
108-86-1	Bromobenzene	0.05	0.2	< 0.2 U
95-49-8	2-Chlorotoluene	0.04	0.2	< 0.2 U
106-43-4	4-Chlorotoluene	0.07	0.2	< 0.2 U
98-06-6	tert-Butylbenzene	0.06	0.2	< 0.2 U
135-98-8	sec-Butylbenzene	0.08	0.2	< 0.2 U
99-87-6	4-Isopropyltoluene	0.08	0.2	< 0.2 U
104-51-8	n-Butylbenzene	0.11	0.2	< 0.2 U
120-82-1	1,2,4-Trichlorobenzene	0.10	0.5	< 0.5 U
91-20-3	Naphthalene	0.07	0.5	< 0.5 U
87-61-6	1,2,3-Trichlorobenzene	0.09	0.5	< 0.5 U

Reported in $\mu\text{g/L}$ (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	93.2%
d8-Toluene	99.5%
Bromo-fluorobenzene	94.7%
d4-1,2-Dichlorobenzene	101%

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

ORGANICS ANALYSIS DATA SHEET

Volatile s by Purge & Trap GC/MS-Method SW8260C

Page 1 of 2

**ANALYTICAL
RESOURCES
INCORPORATED**


Sample ID: FRP-092011-005

SAMPLE

Lab Sample ID: TN19E

LIMS ID: 11-20549

Matrix: Water

Data Release Authorized: *MM*

Reported: 10/03/11

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: 09/20/11

Date Received: 09/20/11

Instrument/Analyst: NT5/PAB

Date Analyzed: 09/21/11 19:15

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.10	0.5	< 0.5 U
74-83-9	Bromomethane	0.04	1.0	< 1.0 U
75-01-4	Vinyl Chloride	0.08	0.2	< 0.2 U
75-00-3	Chloroethane	0.15	0.2	< 0.2 U
75-09-2	Methylene Chloride	0.39	0.5	0.8
67-64-1	Acetone	0.72	5.0	6.1
75-15-0	Carbon Disulfide	0.09	0.2	< 0.2 U
75-35-4	1,1-Dichloroethene	0.09	0.2	< 0.2 U
75-34-3	1,1-Dichloroethane	0.05	0.2	< 0.2 U
156-60-5	trans-1,2-Dichloroethene	0.08	0.2	< 0.2 U
156-59-2	cis-1,2-Dichloroethene	0.10	0.2	< 0.2 U
67-66-3	Chloroform	0.08	0.2	2.0
107-06-2	1,2-Dichloroethane	0.08	0.2	< 0.2 U
78-93-3	2-Butanone	0.81	5.0	< 5.0 U
71-55-6	1,1,1-Trichloroethane	0.09	0.2	< 0.2 U
56-23-5	Carbon Tetrachloride	0.08	0.2	< 0.2 U
108-05-4	Vinyl Acetate	0.07	1.0	< 1.0 U
75-27-4	Bromodichloromethane	0.05	0.2	< 0.2 U
78-87-5	1,2-Dichloropropane	0.09	0.2	< 0.2 U
10061-01-5	cis-1,3-Dichloropropene	0.06	0.2	< 0.2 U
79-01-6	Trichloroethene	0.08	0.2	< 0.2 U
124-48-1	Dibromochloromethane	0.09	0.2	< 0.2 U
79-00-5	1,1,2-Trichloroethane	0.04	0.2	< 0.2 U
71-43-2	Benzene	0.06	0.2	< 0.2 U
10061-02-6	trans-1,3-Dichloropropene	0.06	0.2	< 0.2 U
110-75-8	2-Chloroethylvinylether	0.09	1.0	< 1.0 U
75-25-2	Bromoform	0.07	0.2	< 0.2 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	0.38	5.0	< 5.0 U
591-78-6	2-Hexanone	0.31	5.0	< 5.0 U
127-18-4	Tetrachloroethene	0.09	0.2	< 0.2 U
79-34-5	1,1,2,2-Tetrachloroethane	0.07	0.2	< 0.2 U
108-88-3	Toluene	0.06	0.2	< 0.2 U
108-90-7	Chlorobenzene	0.04	0.2	< 0.2 U
100-41-4	Ethylbenzene	0.09	0.2	< 0.2 U
100-42-5	Styrene	0.07	0.2	< 0.2 U
75-69-4	Trichlorofluoromethane	0.09	0.2	< 0.2 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoro	0.11	0.2	< 0.2 U
179601-23-1	m,p-Xylene	0.14	0.4	< 0.4 U
95-47-6	o-Xylene	0.06	0.2	< 0.2 U
95-50-1	1,2-Dichlorobenzene	0.06	0.2	< 0.2 U
541-73-1	1,3-Dichlorobenzene	0.04	0.2	< 0.2 U
106-46-7	1,4-Dichlorobenzene	0.06	0.2	< 0.2 U
107-02-8	Acrolein	0.29	5.0	< 5.0 U
74-88-4	Methyl Iodide	0.04	1.0	< 1.0 U
74-96-4	Bromoethane	0.09	0.2	< 0.2 U
107-13-1	Acrylonitrile	0.18	1.0	< 1.0 U
563-58-6	1,1-Dichloropropene	0.09	0.2	< 0.2 U
74-95-3	Dibromomethane	0.08	0.2	< 0.2 U
630-20-6	1,1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
96-12-8	1,2-Dibromo-3-chloropropane	0.21	0.5	< 0.5 U

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2

Sample ID: FRP-092011-005
SAMPLE

Lab Sample ID: TN19E
LIMS ID: 11-20549
Matrix: Water
Date Analyzed: 09/21/11 19:15

QC Report No: TN19-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	0.23	0.5	< 0.5 U
110-57-6	trans-1,4-Dichloro-2-butene	0.24	1.0	< 1.0 U
108-67-8	1,3,5-Trimethylbenzene	0.06	0.2	< 0.2 U
95-63-6	1,2,4-Trimethylbenzene	0.06	0.2	< 0.2 U
87-68-3	Hexachlorobutadiene	0.11	0.5	< 0.5 U
106-93-4	Ethylene Dibromide	0.08	0.2	< 0.2 U
74-97-5	Bromoform	0.07	0.2	< 0.2 U
594-20-7	2,2-Dichloropropane	0.08	0.2	< 0.2 U
142-28-9	1,3-Dichloropropane	0.02	0.2	< 0.2 U
98-82-8	Isopropylbenzene	0.06	0.2	< 0.2 U
103-65-1	n-Propylbenzene	0.08	0.2	< 0.2 U
108-86-1	Bromobenzene	0.05	0.2	< 0.2 U
95-49-8	2-Chlorotoluene	0.04	0.2	< 0.2 U
106-43-4	4-Chlorotoluene	0.07	0.2	< 0.2 U
98-06-6	tert-Butylbenzene	0.06	0.2	< 0.2 U
135-98-8	sec-Butylbenzene	0.08	0.2	< 0.2 U
99-87-6	4-Isopropyltoluene	0.08	0.2	< 0.2 U
104-51-8	n-Butylbenzene	0.11	0.2	< 0.2 U
120-82-1	1,2,4-Trichlorobenzene	0.10	0.5	< 0.5 U
91-20-3	Naphthalene	0.07	0.5	< 0.5 U
87-61-6	1,2,3-Trichlorobenzene	0.09	0.5	< 0.5 U

Reported in µg/L (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	109%
d8-Toluene	101%
Bromofluorobenzene	88.7%
d4-1,2-Dichlorobenzene	99.6%

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

ORGANICS ANALYSIS DATA SHEET

Volatile by Purge & Trap GC/MS-Method SW8260C

Page 1 of 2

Sample ID: Trip Blanks
SAMPLE

Lab Sample ID: TN19F

LIMS ID: 11-20550

Matrix: Water

Data Release Authorized: **WW**

Reported: 10/03/11

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: 09/20/11

Date Received: 09/20/11

Instrument/Analyst: NT5/PAB

Date Analyzed: 09/21/11 19:44

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.10	0.5	< 0.5 U
74-83-9	Bromomethane	0.04	1.0	< 1.0 U
75-01-4	Vinyl Chloride	0.08	0.2	< 0.2 U
75-00-3	Chloroethane	0.15	0.2	< 0.2 U
75-09-2	Methylene Chloride	0.39	0.5	< 0.5 U
67-64-1	Acetone	0.72	5.0	< 5.0 U
75-15-0	Carbon Disulfide	0.09	0.2	< 0.2 U
75-35-4	1,1-Dichloroethene	0.09	0.2	< 0.2 U
75-34-3	1,1-Dichloroethane	0.05	0.2	< 0.2 U
156-60-5	trans-1,2-Dichloroethene	0.08	0.2	< 0.2 U
156-59-2	cis-1,2-Dichloroethene	0.10	0.2	< 0.2 U
67-66-3	Chloroform	0.08	0.2	< 0.2 U
107-06-2	1,2-Dichloroethane	0.08	0.2	< 0.2 U
78-93-3	2-Butanone	0.81	5.0	< 5.0 U
71-55-6	1,1,1-Trichloroethane	0.09	0.2	< 0.2 U
56-23-5	Carbon Tetrachloride	0.08	0.2	< 0.2 U
108-05-4	Vinyl Acetate	0.07	1.0	< 1.0 U
75-27-4	Bromodichloromethane	0.05	0.2	< 0.2 U
78-87-5	1,2-Dichloropropane	0.09	0.2	< 0.2 U
10061-01-5	cis-1,3-Dichloropropene	0.06	0.2	< 0.2 U
79-01-6	Trichloroethene	0.08	0.2	< 0.2 U
124-48-1	Dibromochloromethane	0.09	0.2	< 0.2 U
79-00-5	1,1,2-Trichloroethane	0.04	0.2	< 0.2 U
71-43-2	Benzene	0.06	0.2	< 0.2 U
10061-02-6	trans-1,3-Dichloropropene	0.06	0.2	< 0.2 U
110-75-8	2-Chloroethylvinylether	0.09	1.0	< 1.0 U
75-25-2	Bromoform	0.07	0.2	< 0.2 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	0.38	5.0	< 5.0 U
591-78-6	2-Hexanone	0.31	5.0	< 5.0 U
127-18-4	Tetrachloroethene	0.09	0.2	< 0.2 U
79-34-5	1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
108-88-3	Toluene	0.06	0.2	< 0.2 U
108-90-7	Chlorobenzene	0.04	0.2	< 0.2 U
100-41-4	Ethylbenzene	0.09	0.2	< 0.2 U
100-42-5	Styrene	0.07	0.2	< 0.2 U
75-69-4	Trichlorofluoromethane	0.09	0.2	< 0.2 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoro	0.11	0.2	< 0.2 U
179601-23-1	m,p-Xylene	0.14	0.4	< 0.4 U
95-47-6	o-Xylene	0.06	0.2	< 0.2 U
95-50-1	1,2-Dichlorobenzene	0.06	0.2	< 0.2 U
541-73-1	1,3-Dichlorobenzene	0.04	0.2	< 0.2 U
106-46-7	1,4-Dichlorobenzene	0.06	0.2	< 0.2 U
107-02-8	Acrolein	0.29	5.0	< 5.0 U
74-88-4	Methyl Iodide	0.04	1.0	< 1.0 U
74-96-4	Bromoethane	0.09	0.2	< 0.2 U
107-13-1	Acrylonitrile	0.18	1.0	< 1.0 U
563-58-6	1,1-Dichloropropene	0.09	0.2	< 0.2 U
74-95-3	Dibromomethane	0.08	0.2	< 0.2 U
630-20-6	1,1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
96-12-8	1,2-Dibromo-3-chloropropane	0.21	0.5	< 0.5 U

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2

Sample ID: Trip Blanks
SAMPLE

Lab Sample ID: TN19F
LIMS ID: 11-20550
Matrix: Water
Date Analyzed: 09/21/11 19:44

QC Report No: TN19-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	0.23	0.5	< 0.5 U
110-57-6	trans-1,4-Dichloro-2-butene	0.24	1.0	< 1.0 U
108-67-8	1,3,5-Trimethylbenzene	0.06	0.2	< 0.2 U
95-63-6	1,2,4-Trimethylbenzene	0.06	0.2	< 0.2 U
87-68-3	Hexachlorobutadiene	0.11	0.5	< 0.5 U
106-93-4	Ethylene Dibromide	0.08	0.2	< 0.2 U
74-97-5	Bromo-chloromethane	0.07	0.2	< 0.2 U
594-20-7	2,2-Dichloropropane	0.08	0.2	< 0.2 U
142-28-9	1,3-Dichloropropane	0.02	0.2	< 0.2 U
98-82-8	Isopropylbenzene	0.06	0.2	< 0.2 U
103-65-1	n-Propylbenzene	0.08	0.2	< 0.2 U
108-86-1	Bromobenzene	0.05	0.2	< 0.2 U
95-49-8	2-Chlorotoluene	0.04	0.2	< 0.2 U
106-43-4	4-Chlorotoluene	0.07	0.2	< 0.2 U
98-06-6	tert-Butylbenzene	0.06	0.2	< 0.2 U
135-98-8	sec-Butylbenzene	0.08	0.2	< 0.2 U
99-87-6	4-Isopropyltoluene	0.08	0.2	< 0.2 U
104-51-8	n-Butylbenzene	0.11	0.2	< 0.2 U
120-82-1	1,2,4-Trichlorobenzene	0.10	0.5	< 0.5 U
91-20-3	Naphthalene	0.07	0.5	< 0.5 U
87-61-6	1,2,3-Trichlorobenzene	0.09	0.5	< 0.5 U

Reported in $\mu\text{g/L}$ (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	112%
d8-Toluene	102%
Bromo-fluorobenzene	88.9%
d4-1,2-Dichlorobenzene	103%

2-Chloroethylvinylether is an acid labile compound and may not be recovered from an acid preserved sample.

VOA SURROGATE RECOVERY SUMMARY
Matrix: Water
QC Report No: TN19-AMEC Geomatrix
**Project: FRP 2011 Shoreline Investigation
8769**

ARI ID	Client ID	PV	DCE	TOL	BFB	DCB	TOT OUT
TN19A	FRP-092011-001	10	102%	100%	93.8%	98.1%	0
MB-092611	Method Blank	10	94.0%	99.0%	96.0%	100%	0
LCS-092611	Lab Control	10	93.9%	98.3%	98.1%	101%	0
LCSD-092611	Lab Control Dup	10	93.3%	99.9%	98.3%	99.9%	0
TN19B	FRP-092011-002	10	93.9%	100%	96.3%	102%	0
TN19C	FRP-092011-003	10	93.9%	99.5%	95.7%	102%	0
TN19D	FRP-092011-004	10	93.2%	99.5%	94.7%	101%	0
MB-092111	Method Blank	10	110%	102%	95.8%	101%	0
LCS-092111	Lab Control	10	105%	101%	104%	95.8%	0
LCSD-092111	Lab Control Dup	10	104%	99.7%	100%	95.4%	0
TN19E	FRP-092011-005	10	109%	101%	88.7%	99.6%	0
TN19F	Trip Blanks	10	112%	102%	88.9%	103%	0

LCS/MB LIMITS **QC LIMITS**
SW8260C

(DCE) = d4-1,2-Dichloroethane
 (TOL) = d8-Toluene
 (BFB) = Bromofluorobenzene
 (DCB) = d4-1,2-Dichlorobenzene

80-120 80-120
 80-120 80-120
 80-120 80-120
 80-120 80-120

Prep Method: SW5030B
 Log Number Range: 11-20545 to 11-20550

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C

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**ANALYTICAL
RESOURCES
INCORPORATED**


Sample ID: LCS-092111

LAB CONTROL SAMPLE

Lab Sample ID: LCS-092111

LIMS ID: 11-20549

Matrix: Water

Data Release Authorized: *MW*

Reported: 10/03/11

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: NA

Date Received: NA

Instrument/Analyst LCS: NT5/PAB

Sample Amount LCS: 10.0 mL

LCSD: NT5/PAB

LCSD: 10.0 mL

Date Analyzed LCS: 09/21/11 10:01

Purge Volume LCS: 10.0 mL

LCSD: 09/21/11 10:29

LCSD: 10.0 mL

Analyte	LCS	Spike Added-LCS	LCS Recovery	LCSD	Spike Added-LCSD	LCSD Recovery	RPD
Chloromethane	10.7	10.0	107%	11.2	10.0	112%	4.6%
Bromomethane	9.2	10.0	92.0%	10.6	10.0	106%	14.1%
Vinyl Chloride	8.4	10.0	84.0%	10.0	10.0	100%	17.4%
Chloroethane	9.2	10.0	92.0%	9.9	10.0	99.0%	7.3%
Methylene Chloride	8.7	10.0	87.0%	9.9	10.0	99.0%	12.9%
Acetone	59.4	50.0	119%	51.9	50.0	104%	13.5%
Carbon Disulfide	9.5	10.0	95.0%	9.6	10.0	96.0%	1.0%
1,1-Dichloroethene	9.7	10.0	97.0%	10.6	10.0	106%	8.9%
1,1-Dichloroethane	9.9	10.0	99.0%	10.6	10.0	106%	6.8%
trans-1,2-Dichloroethene	9.9	10.0	99.0%	10.4	10.0	104%	4.9%
cis-1,2-Dichloroethene	9.5	10.0	95.0%	10.6	10.0	106%	10.9%
Chloroform	9.8	10.0	98.0%	10.6	10.0	106%	7.8%
1,2-Dichloroethane	9.8	10.0	98.0%	10.8	10.0	108%	9.7%
2-Butanone	54.6	50.0	109%	53.0	50.0	106%	3.0%
1,1,1-Trichloroethane	10.2	10.0	102%	10.8	10.0	108%	5.7%
Carbon Tetrachloride	10.1	10.0	101%	10.9	10.0	109%	7.6%
Vinyl Acetate	9.6	10.0	96.0%	10.5	10.0	105%	9.0%
Bromodichloromethane	9.7	10.0	97.0%	10.6	10.0	106%	8.9%
1,2-Dichloropropane	9.9	10.0	99.0%	10.6	10.0	106%	6.8%
cis-1,3-Dichloropropene	9.7	10.0	97.0%	10.9	10.0	109%	11.7%
Trichloroethene	9.9	10.0	99.0%	10.7	10.0	107%	7.8%
Dibromochloromethane	10.1	10.0	101%	11.0	10.0	110%	8.5%
1,1,2-Trichloroethane	9.9	10.0	99.0%	10.5	10.0	105%	5.9%
Benzene	9.7	10.0	97.0%	10.6	10.0	106%	8.9%
trans-1,3-Dichloropropene	10.2	10.0	102%	11.1	10.0	111%	8.5%
2-Chloroethylvinylether	10.1	10.0	101%	11.2	10.0	112%	10.3%
Bromoform	9.9	10.0	99.0%	11.2	10.0	112%	12.3%
4-Methyl-2-Pentanone (MIBK)	50.3	50.0	101%	53.9	50.0	108%	6.9%
2-Hexanone	54.4	50.0	109%	55.5	50.0	111%	2.0%
Tetrachloroethene	9.9	10.0	99.0%	10.8	10.0	108%	8.7%
1,1,2-Tetrachloroethane	10.0	10.0	100%	11.5	10.0	115%	14.0%
Toluene	10.2	10.0	102%	11.2	10.0	112%	9.3%
Chlorobenzene	10.0	10.0	100%	11.0	10.0	110%	9.5%
Ethylbenzene	10.4	10.0	104%	11.3	10.0	113%	8.3%
Styrene	10.7	10.0	107%	11.6	10.0	116%	8.1%
Trichlorofluoromethane	10.2	10.0	102%	11.0	10.0	110%	7.5%
1,1,2-Trichloro-1,2,2-trifluoroetha	9.6	10.0	96.0%	10.4	10.0	104%	8.0%
m,p-Xylene	21.2	20.0	106%	23.4	20.0	117%	9.9%
o-Xylene	10.5	10.0	105%	11.8	10.0	118%	11.7%
1,2-Dichlorobenzene	9.3	10.0	93.0%	10.6	10.0	106%	13.1%
1,3-Dichlorobenzene	9.3	10.0	93.0%	10.8	10.0	108%	14.9%
1,4-Dichlorobenzene	9.0	10.0	90.0%	10.6	10.0	106%	16.3%
Acrolein	44.6	50.0	89.2%	46.8	50.0	93.6%	4.8%
Methyl Iodide	9.3	10.0	93.0%	10.3	10.0	103%	10.2%
Bromoethane	9.5	10.0	95.0%	10.4	10.0	104%	9.0%

ORGANICS ANALYSIS DATA SHEET

Volatile by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2

ANALYTICAL
RESOURCES
INCORPORATED

Sample ID: LCS-092111
LAB CONTROL SAMPLE

Lab Sample ID: LCS-092111
LIMS ID: 11-20549
Matrix: WaterQC Report No: TN19-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

Analyte	LCS	Spike Added-LCS	LCS Recovery	LCSD	Spike Added-LCSD	LCSD Recovery	RPD
Acrylonitrile	10.7	10.0	107%	10.1	10.0	101%	5.8%
1,1-Dichloropropene	9.6	10.0	96.0%	10.6	10.0	106%	9.9%
Dibromomethane	9.8	10.0	98.0%	10.3	10.0	103%	5.0%
1,1,1,2-Tetrachloroethane	10.3	10.0	103%	11.0	10.0	110%	6.6%
1,2-Dibromo-3-chloropropane	8.1	10.0	81.0%	9.2	10.0	92.0%	12.7%
1,2,3-Trichloropropane	8.3	10.0	83.0%	9.4	10.0	94.0%	12.4%
trans-1,4-Dichloro-2-butene	9.2	10.0	92.0%	10.5	10.0	105%	13.2%
1,3,5-Trimethylbenzene	10.2	10.0	102%	12.0	10.0	120%	16.2%
1,2,4-Trimethylbenzene	10.1	10.0	101%	11.7	10.0	117%	14.7%
Hexachlorobutadiene	8.2 Q	10.0	82.0%	9.6 Q	10.0	96.0%	15.7%
Ethylene Dibromide	9.7	10.0	97.0%	10.3	10.0	103%	6.0%
Bromochloromethane	10.4	10.0	104%	10.8	10.0	108%	3.8%
2,2-Dichloropropane	10.6	10.0	106%	11.5	10.0	115%	8.1%
1,3-Dichloropropane	10.1	10.0	101%	10.7	10.0	107%	5.8%
Isopropylbenzene	10.3	10.0	103%	12.2	10.0	122%	16.9%
n-Propylbenzene	10.0	10.0	100%	11.6	10.0	116%	14.8%
Bromobenzene	9.5	10.0	95.0%	11.1	10.0	111%	15.5%
2-Chlorotoluene	9.8	10.0	98.0%	11.4	10.0	114%	15.1%
4-Chlorotoluene	9.8	10.0	98.0%	11.4	10.0	114%	15.1%
tert-Butylbenzene	9.9	10.0	99.0%	11.5	10.0	115%	15.0%
sec-Butylbenzene	9.7	10.0	97.0%	11.4	10.0	114%	16.1%
4-Isopropyltoluene	10.2	10.0	102%	12.0	10.0	120%	16.2%
n-Butylbenzene	9.9	10.0	99.0%	11.4	10.0	114%	14.1%
1,2,4-Trichlorobenzene	8.9	10.0	89.0%	10.2	10.0	102%	13.6%
Naphthalene	9.2	10.0	92.0%	10.6	10.0	106%	14.1%
1,2,3-Trichlorobenzene	9.7	10.0	97.0%	12.0	10.0	120%	21.2%

Reported in µg/L (ppb)

RPD calculated using sample concentrations per SW846.

Volatile Surrogate Recovery

	LCS	LCSD
d4-1,2-Dichloroethane	105%	104%
d8-Toluene	101%	99.7%
Bromofluorobenzene	104%	100%
d4-1,2-Dichlorobenzene	95.8%	95.4%

ORGANICS ANALYSIS DATA SHEET

Volatile by Purge & Trap GC/MS-Method SW8260C
Page 1 of 2
**ANALYTICAL
RESOURCES
INCORPORATED**


Sample ID: LCS-092611

LAB CONTROL SAMPLE

Lab Sample ID: LCS-092611

LIMS ID: 11-20546

Matrix: Water

Data Release Authorized: *TMW*

Reported: 10/03/11

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: NA

Date Received: NA

Instrument/Analyst LCS: NT2/PAB

LCSD: NT2/PAB

Date Analyzed LCS: 09/26/11 10:06

LCSD: 09/26/11 10:33

Sample Amount LCS: 10.0 mL

LCSD: 10.0 mL

Purge Volume LCS: 10.0 mL

LCSD: 10.0 mL

Analyte	LCS	Spike Added-LCS	LCS Recovery	LCSD	Spike Added-LCSD	LCSD Recovery	RPD
Chloromethane	8.4	10.0	84.0%	8.9	10.0	89.0%	5.8%
Bromomethane	8.0	10.0	80.0%	8.3	10.0	83.0%	3.7%
Vinyl Chloride	9.7	10.0	97.0%	9.7	10.0	97.0%	0.0%
Chloroethane	8.8	10.0	88.0%	9.1	10.0	91.0%	3.4%
Methylene Chloride	9.2	10.0	92.0%	9.5	10.0	95.0%	3.2%
Acetone	41.3 Q	50.0	82.6%	41.9 Q	50.0	83.8%	1.4%
Carbon Disulfide	9.5	10.0	95.0%	9.6	10.0	96.0%	1.0%
1,1-Dichloroethene	8.8	10.0	88.0%	9.2	10.0	92.0%	4.4%
1,1-Dichloroethane	9.2	10.0	92.0%	9.3	10.0	93.0%	1.1%
trans-1,2-Dichloroethene	9.5	10.0	95.0%	9.5	10.0	95.0%	0.0%
cis-1,2-Dichloroethene	9.4	10.0	94.0%	9.7	10.0	97.0%	3.1%
Chloroform	9.5	10.0	95.0%	9.7	10.0	97.0%	2.1%
1,2-Dichloroethane	9.1	10.0	91.0%	9.4	10.0	94.0%	3.2%
2-Butanone	50.0	50.0	100%	49.8	50.0	99.6%	0.4%
1,1,1-Trichloroethane	9.4	10.0	94.0%	9.6	10.0	96.0%	2.1%
Carbon Tetrachloride	9.7	10.0	97.0%	9.9	10.0	99.0%	2.0%
Vinyl Acetate	9.6	10.0	96.0%	9.8	10.0	98.0%	2.1%
Bromodichloromethane	9.6	10.0	96.0%	9.8	10.0	98.0%	2.1%
1,2-Dichloropropane	9.3	10.0	93.0%	9.5	10.0	95.0%	2.1%
cis-1,3-Dichloropropene	9.6	10.0	96.0%	9.8	10.0	98.0%	2.1%
Trichloroethene	9.7	10.0	97.0%	9.9	10.0	99.0%	2.0%
Dibromochloromethane	10.0	10.0	100%	10.0	10.0	100%	0.0%
1,1,2-Trichloroethane	9.6	10.0	96.0%	9.9	10.0	99.0%	3.1%
Benzene	9.4	10.0	94.0%	9.8	10.0	98.0%	4.2%
trans-1,3-Dichloropropene	9.6	10.0	96.0%	9.6	10.0	96.0%	0.0%
2-Chloroethylvinylether	7.7	10.0	77.0%	8.3	10.0	83.0%	7.5%
Bromoform	10.4	10.0	104%	10.3	10.0	103%	1.0%
4-Methyl-2-Pentanone (MIBK)	49.8	50.0	99.6%	51.4	50.0	103%	3.2%
2-Hexanone	49.1	50.0	98.2%	49.6	50.0	99.2%	1.0%
Tetrachloroethene	9.9	10.0	99.0%	9.9	10.0	99.0%	0.0%
1,1,2,2-Tetrachloroethane	10.0	10.0	100%	9.9	10.0	99.0%	1.0%
Toluene	9.6	10.0	96.0%	9.9	10.0	99.0%	3.1%
Chlorobenzene	9.5	10.0	95.0%	9.7	10.0	97.0%	2.1%
Ethylbenzene	9.5	10.0	95.0%	9.5	10.0	95.0%	0.0%
Styrene	10.0	10.0	100%	10.2	10.0	102%	2.0%
Trichlorofluoromethane	9.7	10.0	97.0%	9.8	10.0	98.0%	1.0%
1,1,2-Trichloro-1,2,2-trifluoroethane	9.6	10.0	96.0%	9.7	10.0	97.0%	1.0%
m,p-Xylene	19.5	20.0	97.5%	19.8	20.0	99.0%	1.5%
o-Xylene	9.6	10.0	96.0%	9.7	10.0	97.0%	1.0%
1,2-Dichlorobenzene	9.6	10.0	96.0%	9.7	10.0	97.0%	1.0%
1,3-Dichlorobenzene	9.7	10.0	97.0%	9.7	10.0	97.0%	0.0%
1,4-Dichlorobenzene	9.6	10.0	96.0%	9.7	10.0	97.0%	1.0%
Acrolein	53.6	50.0	107%	56.1	50.0	112%	4.6%
Methyl Iodide	7.6 Q	10.0	76.0%	8.0 Q	10.0	80.0%	5.1%
Bromoethane	9.5	10.0	95.0%	9.6	10.0	96.0%	1.0%

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2

Sample ID: LCS-092611
LAB CONTROL SAMPLE

Lab Sample ID: LCS-092611
LIMS ID: 11-20546
Matrix: Water

QC Report No: TN19-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

Analyte	LCS	Spike Added-LCS	LCS Recovery	LCSD	Spike Added-LCSD	LCSD Recovery	RPD
Acrylonitrile	9.4	10.0	94.0%	9.7	10.0	97.0%	3.1%
1,1-Dichloropropene	9.6	10.0	96.0%	9.9	10.0	99.0%	3.1%
Dibromomethane	9.4	10.0	94.0%	9.7	10.0	97.0%	3.1%
1,1,1,2-Tetrachloroethane	9.9	10.0	99.0%	9.9	10.0	99.0%	0.0%
1,2-Dibromo-3-chloropropane	9.6	10.0	96.0%	9.4	10.0	94.0%	2.1%
1,2,3-Trichloropropane	9.5	10.0	95.0%	9.6	10.0	96.0%	1.0%
trans-1,4-Dichloro-2-butene	7.8 Q	10.0	78.0%	7.7 Q	10.0	77.0%	1.3%
1,3,5-Trimethylbenzene	9.7	10.0	97.0%	9.7	10.0	97.0%	0.0%
1,2,4-Trimethylbenzene	9.7	10.0	97.0%	9.9	10.0	99.0%	2.0%
Hexachlorobutadiene	9.8	10.0	98.0%	10.3	10.0	103%	5.0%
Ethylene Dibromide	9.6	10.0	96.0%	9.9	10.0	99.0%	3.1%
Bromochloromethane	9.7	10.0	97.0%	9.8	10.0	98.0%	1.0%
2,2-Dichloropropane	9.0	10.0	90.0%	9.0	10.0	90.0%	0.0%
1,3-Dichloropropane	9.5	10.0	95.0%	9.5	10.0	95.0%	0.0%
Isopropylbenzene	9.7	10.0	97.0%	9.7	10.0	97.0%	0.0%
n-Propylbenzene	9.5	10.0	95.0%	9.7	10.0	97.0%	2.1%
Bromobenzene	9.7	10.0	97.0%	9.7	10.0	97.0%	0.0%
2-Chlorotoluene	9.4	10.0	94.0%	9.5	10.0	95.0%	1.1%
4-Chlorotoluene	9.5	10.0	95.0%	9.6	10.0	96.0%	1.0%
tert-Butylbenzene	9.6	10.0	96.0%	9.7	10.0	97.0%	1.0%
sec-Butylbenzene	9.5	10.0	95.0%	9.8	10.0	98.0%	3.1%
4-Isopropyltoluene	9.6	10.0	96.0%	10.0	10.0	100%	4.1%
n-Butylbenzene	9.0	10.0	90.0%	9.7	10.0	97.0%	7.5%
1,2,4-Trichlorobenzene	9.8	10.0	98.0%	9.9	10.0	99.0%	1.0%
Naphthalene	10.2	10.0	102%	10.3	10.0	103%	1.0%
1,2,3-Trichlorobenzene	10.0	10.0	100%	10.2	10.0	102%	2.0%

Reported in $\mu\text{g/L}$ (ppb)

RPD calculated using sample concentrations per SW846.

Volatile Surrogate Recovery

	LCS	LCSD
d4-1,2-Dichloroethane	93.9%	93.3%
d8-Toluene	98.3%	99.9%
Bromofluorobenzene	98.1%	98.3%
d4-1,2-Dichlorobenzene	101%	99.9%

4A
VOLATILE METHOD BLANK SUMMARY

Method Blank ID.

MB0921

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONME

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVE

Lab File ID: MB0921

Lab Sample ID: MB0921

Date Analyzed: 09/21/11

Time Analyzed: 1057

Instrument ID: NT5

Heated Purge: (Y/N) N

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01 LCS0921	LCS0921	LCS0921	1001
02 LCS0921	LCS0921	LCS0921A	1029
03 FRP-091911-0	TN00B	TN00B	1433
04 FRP-091911-0	TN00C	TN00C	1501
05 FRP-091911-0	TN00D	TN00D	1529
06 FRP-091911-0	TN00E	TN00E	1557
07 FRP-091911-0	TN00F	TN00F	1626
08 TRIP BLANKS	TN00G	TN00G	1654
09 FRP-092011-0	TN19A	TN19A	1722
10 FRP-092011-0	TN19E	TN19E	1915
11 TRIP BLANKS	TN19F	TN19F	1944
12			
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COMMENTS:

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 1 of 2

ANALYTICAL
RESOURCES
INCORPORATED

Sample ID: MB-092111

METHOD BLANK

Lab Sample ID: MB-092111
LIMS ID: 11-20549
Matrix: Water
Data Release Authorized: *MW*
Reported: 10/03/11

QC Report No: TN19-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769
Date Sampled: NA
Date Received: NA

Instrument/Analyst: NT5/PAB
Date Analyzed: 09/21/11 10:57

Sample Amount: 10.0 mL
Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.10	0.5	< 0.5 U
74-83-9	Bromomethane	0.04	1.0	< 1.0 U
75-01-4	Vinyl Chloride	0.08	0.2	< 0.2 U
75-00-3	Chloroethane	0.15	0.2	< 0.2 U
75-09-2	Methylene Chloride	0.39	0.5	< 0.5 U
67-64-1	Acetone	0.72	5.0	< 5.0 U
75-15-0	Carbon Disulfide	0.09	0.2	< 0.2 U
75-35-4	1,1-Dichloroethene	0.09	0.2	< 0.2 U
75-34-3	1,1-Dichloroethane	0.05	0.2	< 0.2 U
156-60-5	trans-1,2-Dichloroethene	0.08	0.2	< 0.2 U
156-59-2	cis-1,2-Dichloroethene	0.10	0.2	< 0.2 U
67-66-3	Chloroform	0.08	0.2	< 0.2 U
107-06-2	1,2-Dichloroethane	0.08	0.2	< 0.2 U
78-93-3	2-Butanone	0.81	5.0	< 5.0 U
71-55-6	1,1,1-Trichloroethane	0.09	0.2	< 0.2 U
56-23-5	Carbon Tetrachloride	0.08	0.2	< 0.2 U
108-05-4	Vinyl Acetate	0.07	1.0	< 1.0 U
75-27-4	Bromodichloromethane	0.05	0.2	< 0.2 U
78-87-5	1,2-Dichloropropane	0.09	0.2	< 0.2 U
10061-01-5	cis-1,3-Dichloropropene	0.06	0.2	< 0.2 U
79-01-6	Trichloroethene	0.08	0.2	< 0.2 U
124-48-1	Dibromochloromethane	0.09	0.2	< 0.2 U
79-00-5	1,1,2-Trichloroethane	0.04	0.2	< 0.2 U
71-43-2	Benzene	0.06	0.2	< 0.2 U
10061-02-6	trans-1,3-Dichloropropene	0.06	0.2	< 0.2 U
110-75-8	2-Chloroethylvinylether	0.09	1.0	< 1.0 U
75-25-2	Bromoform	0.07	0.2	< 0.2 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	0.38	5.0	< 5.0 U
591-78-6	2-Hexanone	0.31	5.0	< 5.0 U
127-18-4	Tetrachloroethene	0.09	0.2	< 0.2 U
79-34-5	1,1,2,2-Tetrachloroethane	0.07	0.2	< 0.2 U
108-88-3	Toluene	0.06	0.2	< 0.2 U
108-90-7	Chlorobenzene	0.04	0.2	< 0.2 U
100-41-4	Ethylbenzene	0.09	0.2	< 0.2 U
100-42-5	Styrene	0.07	0.2	< 0.2 U
75-69-4	Trichlorofluoromethane	0.09	0.2	< 0.2 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoro	0.11	0.2	< 0.2 U
179601-23-1	m,p-Xylene	0.14	0.4	< 0.4 U
95-47-6	o-Xylene	0.06	0.2	< 0.2 U
95-50-1	1,2-Dichlorobenzene	0.06	0.2	< 0.2 U
541-73-1	1,3-Dichlorobenzene	0.04	0.2	< 0.2 U
106-46-7	1,4-Dichlorobenzene	0.06	0.2	< 0.2 U
107-02-8	Acrolein	0.29	5.0	< 5.0 U
74-88-4	Methyl Iodide	0.04	1.0	< 1.0 U
74-96-4	Bromoethane	0.09	0.2	< 0.2 U
107-13-1	Acrylonitrile	0.18	1.0	< 1.0 U
563-58-6	1,1-Dichloropropene	0.09	0.2	< 0.2 U
74-95-3	Dibromomethane	0.08	0.2	< 0.2 U
630-20-6	1,1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
96-12-8	1,2-Dibromo-3-chloropropane	0.21	0.5	< 0.5 U

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
Page 2 of 2

Sample ID: MB-092111

METHOD BLANK

Lab Sample ID: MB-092111

LIMS ID: 11-20549

Matrix: Water

Date Analyzed: 09/21/11 10:57

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	0.23	0.5	< 0.5 U
110-57-6	trans-1,4-Dichloro-2-butene	0.24	1.0	< 1.0 U
108-67-8	1,3,5-Trimethylbenzene	0.06	0.2	< 0.2 U
95-63-6	1,2,4-Trimethylbenzene	0.06	0.2	< 0.2 U
87-68-3	Hexachlorobutadiene	0.11	0.5	< 0.5 U
106-93-4	Ethylene Dibromide	0.08	0.2	< 0.2 U
74-97-5	Bromochloromethane	0.07	0.2	< 0.2 U
594-20-7	2,2-Dichloropropane	0.08	0.2	< 0.2 U
142-28-9	1,3-Dichloropropane	0.02	0.2	< 0.2 U
98-82-8	Isopropylbenzene	0.06	0.2	< 0.2 U
103-65-1	n-Propylbenzene	0.08	0.2	< 0.2 U
108-86-1	Bromobenzene	0.05	0.2	< 0.2 U
95-49-8	2-Chlorotoluene	0.04	0.2	< 0.2 U
106-43-4	4-Chlorotoluene	0.07	0.2	< 0.2 U
98-06-6	tert-Butylbenzene	0.06	0.2	< 0.2 U
135-98-8	sec-Butylbenzene	0.08	0.2	< 0.2 U
99-87-6	4-Isopropyltoluene	0.08	0.2	< 0.2 U
104-51-8	n-Butylbenzene	0.11	0.2	< 0.2 U
120-82-1	1,2,4-Trichlorobenzene	0.10	0.5	< 0.5 U
91-20-3	Naphthalene	0.07	0.5	< 0.5 U
87-61-6	1,2,3-Trichlorobenzene	0.09	0.5	< 0.5 U

Reported in $\mu\text{g/L}$ (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	110%
d8-Toluene	102%
Bromoform	95.8%
d4-1,2-Dichlorobenzene	101%

4A
VOLATILE METHOD BLANK SUMMARY

Method Blank ID.

Lab Name: ANALYTICAL RESOURCES INC

ARI Job No: TN42

Lab File ID: MB0926

Date Analyzed: 09/26/11

Instrument ID: NT2

	MB0926
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Client: AMEC EARTH AND ENVIRONME

Project: 2011 FRP SHORELINE INVE

Lab Sample ID: MB0926

Time Analyzed: 1100

Heated Purge: (Y/N) N

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	TIME ANALYZED
01 LCS0926	LCS0926	LCS0926	1006
02 LCS0926	LCS0926	LCS0926A	1033
03 TRIP BLANKS	TN42G	TN42G2	1228
04 FRP-091911-0	TN00B	TN00B2	1255
05 FRP-091911-0	TN00C	TN00C2	1322
06 FRP-091911-0	TN00D	TN00D2	1349
07 FRP-091911-0	TN00E	TN00E2	1415
08 FRP-092011-0	TN19B	TN19B2	1442
09 FRP-092011-0	TN19C	TN19C2	1509
10 FRP-092011-0	TN19D	TN19D2	1536
11 FRP-092111-0	TN42A	TN42A2	1603
12 FRP-092111-0	TN42B	TN42B2	1630
13 FRP-092111-0	TN42C	TN42C2	1657
14 FRP-092111-0	TN42D	TN42D2	1723
15 FRP-092111-0	TN42E	TN42E2	1750
16 FRP-092111-0	TN42F	TN42F2	1817
17 FRP-091911-0	TN00CMS	TN00CMS	1843
18 FRP-091911-0	TN00CMSP	TN00CMSP	1910
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COMMENTS:

ORGANICS ANALYSIS DATA SHEET

Volatile by Purge & Trap GC/MS-Method SW8260C
Page 1 of 2

**ANALYTICAL
RESOURCES
INCORPORATED**

Sample ID: MB-092611

METHOD BLANK

Lab Sample ID: MB-092611

LIMS ID: 11-20546

Matrix: Water

Data Release Authorized: 

Reported: 10/03/11

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: NA

Date Received: NA

Instrument/Analyst: NT2/PAB

Date Analyzed: 09/26/11 11:00

Sample Amount: 10.0 mL

Purge Volume: 10.0 mL

CAS Number	Analyte	MDL	RL	Result
74-87-3	Chloromethane	0.10	0.5	< 0.5 U
74-83-9	Bromomethane	0.04	1.0	< 1.0 U
75-01-4	Vinyl Chloride	0.08	0.2	< 0.2 U
75-00-3	Chloroethane	0.15	0.2	< 0.2 U
75-09-2	Methylene Chloride	0.39	0.5	< 0.5 U
67-64-1	Acetone	0.72	5.0	< 5.0 U
75-15-0	Carbon Disulfide	0.09	0.2	< 0.2 U
75-35-4	1,1-Dichloroethene	0.09	0.2	< 0.2 U
75-34-3	1,1-Dichloroethane	0.05	0.2	< 0.2 U
156-60-5	trans-1,2-Dichloroethene	0.08	0.2	< 0.2 U
156-59-2	cis-1,2-Dichloroethene	0.10	0.2	< 0.2 U
67-66-3	Chloroform	0.08	0.2	< 0.2 U
107-06-2	1,2-Dichloroethane	0.08	0.2	< 0.2 U
78-93-3	2-Butanone	0.81	5.0	< 5.0 U
71-55-6	1,1,1-Trichloroethane	0.09	0.2	< 0.2 U
56-23-5	Carbon Tetrachloride	0.08	0.2	< 0.2 U
108-05-4	Vinyl Acetate	0.07	1.0	< 1.0 U
75-27-4	Bromodichloromethane	0.05	0.2	< 0.2 U
78-87-5	1,2-Dichloropropane	0.09	0.2	< 0.2 U
10061-01-5	cis-1,3-Dichloropropene	0.06	0.2	< 0.2 U
79-01-6	Trichloroethene	0.08	0.2	< 0.2 U
124-48-1	Dibromochloromethane	0.09	0.2	< 0.2 U
79-00-5	1,1,2-Trichloroethane	0.04	0.2	< 0.2 U
71-43-2	Benzene	0.06	0.2	< 0.2 U
10061-02-6	trans-1,3-Dichloropropene	0.06	0.2	< 0.2 U
110-75-8	2-Chloroethylvinylether	0.09	1.0	< 1.0 U
75-25-2	Bromoform	0.07	0.2	< 0.2 U
108-10-1	4-Methyl-2-Pentanone (MIBK)	0.38	5.0	< 5.0 U
591-78-6	2-Hexanone	0.31	5.0	< 5.0 U
127-18-4	Tetrachloroethene	0.09	0.2	< 0.2 U
79-34-5	1,1,2,2-Tetrachloroethane	0.07	0.2	< 0.2 U
108-88-3	Toluene	0.06	0.2	< 0.2 U
108-90-7	Chlorobenzene	0.04	0.2	< 0.2 U
100-41-4	Ethylbenzene	0.09	0.2	< 0.2 U
100-42-5	Styrene	0.07	0.2	< 0.2 U
75-69-4	Trichlorofluoromethane	0.09	0.2	< 0.2 U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoro	0.11	0.2	< 0.2 U
179601-23-1	m,p-Xylene	0.14	0.4	< 0.4 U
95-47-6	o-Xylene	0.06	0.2	< 0.2 U
95-50-1	1,2-Dichlorobenzene	0.06	0.2	< 0.2 U
541-73-1	1,3-Dichlorobenzene	0.04	0.2	< 0.2 U
106-46-7	1,4-Dichlorobenzene	0.06	0.2	< 0.2 U
107-02-8	Acrolein	0.29	5.0	< 5.0 U
74-88-4	Methyl Iodide	0.04	1.0	< 1.0 U
74-96-4	Bromoethane	0.09	0.2	< 0.2 U
107-13-1	Acrylonitrile	0.18	1.0	< 1.0 U
563-58-6	1,1-Dichloropropene	0.09	0.2	< 0.2 U
74-95-3	Dibromomethane	0.08	0.2	< 0.2 U
630-20-6	1,1,1,2-Tetrachloroethane	0.07	0.2	< 0.2 U
96-12-8	1,2-Dibromo-3-chloropropane	0.21	0.5	< 0.5 U

ORGANICS ANALYSIS DATA SHEET

Volatiles by Purge & Trap GC/MS-Method SW8260C
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Sample ID: MB-092611

METHOD BLANK

Lab Sample ID: MB-092611
LIMS ID: 11-20546
Matrix: Water
Date Analyzed: 09/26/11 11:00QC Report No: TN19-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

CAS Number	Analyte	MDL	RL	Result
96-18-4	1,2,3-Trichloropropane	0.23	0.5	< 0.5 U
110-57-6	trans-1,4-Dichloro-2-butene	0.24	1.0	< 1.0 U
108-67-8	1,3,5-Trimethylbenzene	0.06	0.2	< 0.2 U
95-63-6	1,2,4-Trimethylbenzene	0.06	0.2	< 0.2 U
87-68-3	Hexachlorobutadiene	0.11	0.5	< 0.5 U
106-93-4	Ethylene Dibromide	0.08	0.2	< 0.2 U
74-97-5	Bromo-chloromethane	0.07	0.2	< 0.2 U
594-20-7	2,2-Dichloropropane	0.08	0.2	< 0.2 U
142-28-9	1,3-Dichloropropane	0.02	0.2	< 0.2 U
98-82-8	Isopropylbenzene	0.06	0.2	< 0.2 U
103-65-1	n-Propylbenzene	0.08	0.2	< 0.2 U
108-86-1	Bromobenzene	0.05	0.2	< 0.2 U
95-49-8	2-Chlorotoluene	0.04	0.2	< 0.2 U
106-43-4	4-Chlorotoluene	0.07	0.2	< 0.2 U
98-06-6	tert-Butylbenzene	0.06	0.2	< 0.2 U
135-98-8	sec-Butylbenzene	0.08	0.2	< 0.2 U
99-87-6	4-Isopropyltoluene	0.08	0.2	< 0.2 U
104-51-8	n-Butylbenzene	0.11	0.2	< 0.2 U
120-82-1	1,2,4-Trichlorobenzene	0.10	0.5	< 0.5 U
91-20-3	Naphthalene	0.07	0.5	< 0.5 U
87-61-6	1,2,3-Trichlorobenzene	0.09	0.5	< 0.5 U

Reported in µg/L (ppb)

Volatile Surrogate Recovery

d4-1,2-Dichloroethane	94.0%
d8-Toluene	99.0%
Bromo-fluorobenzene	96.0%
d4-1,2-Dichlorobenzene	100%

5A
VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: ANALYTICAL RESOURCES INC Contract: AMEC EARTH AND ENVIRONMENTAL
 Lab Code: ARI Case No.: 2011 FRP SHORELINE INVESTIGATION SDG No.: TN42
 Lab File ID: BFB0727 BFB Injection Date: 07/27/11
 Instrument ID: NT5 BFB Injection Time: 1130
 GC Column: RTXVMS ID: 0.18 (mm) Heated Purge: (Y/N) N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	8.0 - 40.0% of mass 95	16.1
75	30.0 - 66.0% of mass 95	55.5
95	Base Peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	7.0
173	Less than 2.0% of mass 174	0.7 (1.0)1
174	50.0 - 101.0% of mass 95	72.0
175	4.0 - 9.0% of mass 174	6.0 (8.4)1
176	93.0 - 101.0% of mass 174	72.1 (100.1)1
177	5.0 - 9.0% of mass 176	4.7 (6.5)2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	VSTD60	IC0727	6000727	07/27/11	1220
02	VSTD40	IC0727	4000727	07/27/11	1248
03	VSTD20	IC0727	2000727	07/27/11	1316
04	VSTD10	IC0727	1000727	07/27/11	1345
05	VSTD4	IC0727	0400727	07/27/11	1413
06	VSTD1	IC0727	0100727	07/27/11	1441
07	VSTD0.5	IC0727	0050727	07/27/11	1510
08	VSTD0.2	IC0727	0020727	07/27/11	1538
09					
10					
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5A
VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: ANALYTICAL RESOURCES INC Contract: AMEC EARTH AND ENVIRONMENTAL
 Lab Code: ARI Case No.: 2011 FRP SHORELINE INVESTIGATION SDG No.: TN42
 Lab File ID: BFB0921 BFB Injection Date: 09/21/11
 Instrument ID: NT5 BFB Injection Time: 0812
 GC Column: RTXVMS ID: 0.18 (mm) Heated Purge: (Y/N) N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	8.0 - 40.0% of mass 95	18.7
75	30.0 - 66.0% of mass 95	55.0
95	Base Peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	5.9
173	Less than 2.0% of mass 174	0.2 (0.2)1
174	50.0 - 101.0% of mass 95	73.7
175	4.0 - 9.0% of mass 174	5.4 (7.3)1
176	93.0 - 101.0% of mass 174	73.3 (99.4)1
177	5.0 - 9.0% of mass 176	4.4 (5.9)2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	VSTD10	CC0921	1000921	09/21/11	0855
02	LCS0921	LCS0921	LCS0921	09/21/11	1001
03	LCS0921	LCS0921	LCS0921A	09/21/11	1029
04	MB0921	MB0921	MB0921	09/21/11	1057
05	FRP-091911-002	TN00B	TN00B	09/21/11	1433
06	FRP-091911-003	TN00C	TN00C	09/21/11	1501
07	FRP-091911-004	TN00D	TN00D	09/21/11	1529
08	FRP-091911-005	TN00E	TN00E	09/21/11	1557
09	FRP-091911-006	TN00F	TN00F	09/21/11	1626
10	TRIP BLANKS	TN00G	TN00G	09/21/11	1654
11	FRP-092011-001	TN19A	TN19A	09/21/11	1722
12	FRP-092011-005	TN19E	TN19E	09/21/11	1915
13	TRIP BLANKS	TN19F	TN19F	09/21/11	1944
14					
15					
16					
17					
18					
19					
20					
21					
22					

5A
 VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
 BROMOFLUOROBENZENE (BFB)

Lab Name: ANALYTICAL RESOURCES INC Contract: AMEC EARTH AND ENVIRONMENTAL
 Lab Code: ARI Case No.: 2011 FRP SHORELINE INVESTIGATION SDG No.: TN42
 Lab File ID: BFB0921 BFB Injection Date: 09/21/11
 Instrument ID: NT2 BFB Injection Time: 1241
 GC Column: RTXVMS ID: 0.18 (mm) Heated Purge: (Y/N) N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	8.0 - 40.0% of mass 95	18.7
75	30.0 - 66.0% of mass 95	53.0
95	Base Peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	7.1
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	50.0 - 101.0% of mass 95	79.2
175	4.0 - 9.0% of mass 174	5.7 (7.2)1
176	93.0 - 101.0% of mass 174	79.4 (100.3)1
177	5.0 - 9.0% of mass 176	5.5 (7.0)2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01 VSTD0.2	VSTD0.2	00_20921	09/21/11	1338
02 VSTD0.5	VSTD0.5	00_50921	09/21/11	1405
03 VSTD01	VSTD01	01_00921	09/21/11	1432
04 VSTD02	VSTD02	02_00921	09/21/11	1459
05 VSTD10	VSTD10	10_00921	09/21/11	1526
06 VSTD20	VSTD20	20_00921	09/21/11	1552
07 VSTD40	VSTD40	40_00921	09/21/11	1619
08 VSTD60	VSTD60	60_00921	09/21/11	1646
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5A
VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: ANALYTICAL RESOURCES INC Contract: AMEC EARTH AND ENVIRONMENTAL
 Lab Code: ARI Case No.: 2011 FRP SHORELINE INVESTIGATION SDG No.: TN42
 Lab File ID: BFB0926 BFB Injection Date: 09/26/11
 Instrument ID: NT2 BFB Injection Time: 0901
 GC Column: RTXVMS ID: 0.18 (mm) Heated Purge: (Y/N) N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	8.0 - 40.0% of mass 95	17.4
75	30.0 - 66.0% of mass 95	51.2
95	Base Peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	7.2
173	Less than 2.0% of mass 174	0.1 (0.1)1
174	50.0 - 101.0% of mass 95	82.2
175	4.0 - 9.0% of mass 174	5.7 (6.9)1
176	93.0 - 101.0% of mass 174	81.9 (99.6)1
177	5.0 - 9.0% of mass 176	5.5 (6.7)2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

	EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	CC0926	CC0926	CC0926	09/26/11	0939
02	LCS0926	LCS0926	LCS0926	09/26/11	1006
03	LCS0926	LCS0926	LCS0926A	09/26/11	1033
04	MB0926	MB0926	MB0926	09/26/11	1100
05	TRIP BLANKS	TN42G	TN42G2	09/26/11	1228
06	FRP-091911-002	TN00B	TN00B2	09/26/11	1255
07	FRP-091911-003	TN00C	TN00C2	09/26/11	1322
08	FRP-091911-004	TN00D	TN00D2	09/26/11	1349
09	FRP-091911-005	TN00E	TN00E2	09/26/11	1415
10	FRP-092011-002	TN19B	TN19B2	09/26/11	1442
11	FRP-092011-003	TN19C	TN19C2	09/26/11	1509
12	FRP-092011-004	TN19D	TN19D2	09/26/11	1536
13	FRP-092111-001	TN42A	TN42A2	09/26/11	1603
14	FRP-092111-002	TN42B	TN42B2	09/26/11	1630
15	FRP-092111-003	TN42C	TN42C2	09/26/11	1657
16	FRP-092111-004	TN42D	TN42D2	09/26/11	1723
17	FRP-092111-005	TN42E	TN42E2	09/26/11	1750
18	FRP-092111-006	TN42F	TN42F2	09/26/11	1817
19	FRP-091911-003MS	TN00CMS	TN00CMS	09/26/11	1843
20	FRP-091911-003MS	TN00CMSD	TN00CMSD	09/26/11	1910
21					
22					

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

LAB FILE ID: RF0.2: 0020727 RF0.5: 0050727 RF1: 0100727
 RF4: 0400727 RF10: 1000727

COMPOUND	RF0.2	RF0.5	RF1	RF4	RF10
Chloromethane		0.445	0.457	0.360	0.340
Vinyl Chloride	0.387	0.360	0.422	0.419	0.431
Bromomethane	0.197	0.224	0.234	0.250	0.262
Chloroethane	0.276	0.227	0.280	0.231	0.255
Trichlorofluoromethane	0.957	0.998	1.258	1.074	1.068
Acrolein		0.042	0.034	0.033	0.034
112Trichloro122Trifluoroetha	0.463	0.450	0.550	0.466	0.445
Acetone			0.061	0.057	0.063
1,1-Dichloroethene	0.352	0.371	0.490	0.397	0.396
Bromoethane	0.262	0.252	0.278	0.248	0.257
Iodomethane		0.564	0.749	0.624	0.606
Methylene Chloride	0.348	0.486	0.452	0.387	0.396
Acrylonitrile			0.089	0.069	0.060
Carbon Disulfide	1.156	1.041	1.184	1.061	1.118
Trans-1,2-Dichloroethene	0.430	0.490	0.510	0.467	0.460
Vinyl Acetate		0.413	0.474	0.465	0.461
1,1-Dichloroethane	0.679	0.751	0.805	0.733	0.736
2-Butanone		0.078	0.097	0.087	0.083
2,2-Dichloropropane	0.862	0.880	1.027	0.962	0.990
Cis-1,2-Dichloroethene	0.446	0.434	0.510	0.463	0.474
Chloroform	1.037	0.831	1.051	0.976	0.911
Bromochloromethane	0.185	0.222	0.241	0.232	0.214
1,1,1-Trichloroethane	0.935	1.018	1.168	1.118	1.080
1,1-Dichloropropene	0.495	0.433	0.524	0.489	0.504
Carbon Tetrachloride	0.648	0.697	0.815	0.779	0.777
1,2-Dichloroethane	0.571	0.510	0.641	0.592	0.574
Benzene	1.219	1.141	1.347	1.264	1.263
Trichloroethene	0.359	0.361	0.405	0.402	0.397
1,2-Dichloropropane	0.197	0.242	0.301	0.233	0.240
Bromodichloromethane	0.587	0.499	0.577	0.564	0.540
Dibromomethane	0.172	0.180	0.213	0.200	0.194
2-Chloroethyl Vinyl Ether		0.094	0.116	0.115	0.124
4-Methyl-2-Pentanone		0.058	0.064	0.063	0.062
Cis 1,3-dichloropropene	0.466	0.524	0.587	0.555	0.548
Toluene	0.906	0.806	0.927	0.902	0.904
Trans 1,3-Dichloropropene	0.529	0.478	0.594	0.558	0.556
2-Hexanone		0.099	0.124	0.114	0.115

FORM VI VOA

TN19 : 000052

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

LAB FILE ID: RF0.2: 0020727 RF0.5: 0050727 RF1: 0100727
 RF4: 0400727 RF10: 1000727

COMPOUND	RF0.2	RF0.5	RF1	RF4	RF10
1,1,2-Trichloroethane	0.227	0.237	0.281	0.249	0.239
1,3-Dichloropropane	0.471	0.429	0.523	0.470	0.439
Tetrachloroethene	0.510	0.433	0.513	0.467	0.454
Chlorodibromomethane	0.381	0.358	0.464	0.425	0.420
1,2-Dibromoethane	0.280	0.234	0.279	0.264	0.262
Chlorobenzene	1.158	1.115	1.204	1.137	1.105
Ethyl Benzene	0.584	0.562	0.651	0.606	0.613
1,1,1,2-Tetrachloroethane	0.450	0.454	0.514	0.471	0.461
m,p-xylene	0.719	0.700	0.823	0.767	0.747
o-Xylene	0.592	0.609	0.769	0.695	0.714
Styrene	0.995	1.042	1.262	1.183	1.210
Bromoform	0.486	0.394	0.515	0.437	0.411
1,1,2,2-Tetrachloroethane	0.675	0.420	0.492	0.469	0.413
1,2,3-Trichloropropane	0.264	0.209	0.283	0.207	0.191
Trans-1,4-Dichloro 2-Butene			0.229	0.236	0.207
N-Propyl Benzene	3.544	3.127	3.884	3.502	3.307
Bromobenzene	0.950	0.906	1.048	0.900	0.835
Isopropyl Benzene	2.811	2.776	3.485	3.289	3.092
2-Chloro Toluene	2.351	2.305	2.711	2.443	2.307
4-Chloro Toluene	2.961	2.400	2.912	2.656	2.526
T-Butyl Benzene	2.160	2.101	2.624	2.386	2.279
1,3,5-Trimethyl Benzene	2.657	2.500	3.198	2.913	2.747
1,2,4-Trimethylbenzene	2.740	2.526	3.143	2.980	2.852
S-Butyl Benzene	3.035	2.770	3.343	2.996	2.938
4-Isopropyl Toluene	2.488	2.252	2.859	2.751	2.701
1,3-Dichlorobenzene	1.678	1.489	1.999	1.701	1.598
1,4-Dichlorobenzene	2.105	1.802	1.971	1.734	1.649
N-Butyl Benzene	1.928	1.951	2.329	2.148	2.111
1,2-Dichlorobenzene	1.733	1.498	1.780	1.579	1.456
1,2-Dibromo 3-Chloropropane	0.104	0.177	0.173	0.133	0.120
1,2,4-Trichlorobenzene	0.910	0.873	1.027	0.890	0.833
Hexachloro 1,3-Butadiene		0.550	0.577	0.472	0.444
Naphthalene	1.426	1.383	1.643	1.510	1.454
1,2,3-Trichlorobenzene	0.750	0.612	0.758	0.638	0.546
Dichlorodifluoromethane	0.511	0.578	0.594	0.616	0.646
Methyl tert butyl ether	1.258	1.246	1.428	1.319	1.291

FORM VI VOA

TN19 : 000053

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

LAB FILE ID: RF0.2: 0020727 RF0.5: 0050727 RF1: 0100727
RF4: 0400727 RF10: 1000727

COMPOUND	RF0.2	RF0.5	RF1	RF4	RF10
d4-1,2-Dichloroethane	0.541	0.588	0.595	0.598	0.570
d8-Toluene	1.120	1.100	1.111	1.136	1.126
4-Bromofluorobenzene	0.554	0.579	0.575	0.589	0.596
d4-1,2-Dichlorobenzene	0.861	0.875	0.902	0.882	0.882
Dibromofluoromethane	0.399	0.403	0.423	0.423	0.426

FORM VI VOA

TN19 : 00054

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

LAB FILE ID: RF20: 2000727

RF40: 4000727

RF60: 6000727

COMPOUND	RF20	RF40	RF60
Chloromethane	0.307	0.274	0.259
Vinyl Chloride	0.382	0.362	0.311
Bromomethane	0.267	0.276	0.256
Chloroethane	0.215	0.210	0.203
Trichlorofluoromethane	0.942	0.762	0.719
Acrolein	0.031	0.029	0.028
112Trichloro122Trifluoroetha	0.409	0.391	0.374
Acetone	0.049	0.054	0.053
1,1-Dichloroethene	0.359	0.339	0.314
Bromoethane	0.233	0.207	0.180
Iodomethane	0.548	0.505	0.418
Methylene Chloride	0.356	0.322	0.283
Acrylonitrile	0.060	0.059	0.058
Carbon Disulfide	0.995	0.897	0.942
Trans-1,2-Dichloroethene	0.437	0.405	0.414
Vinyl Acetate	0.456	0.463	0.423
1,1-Dichloroethane	0.690	0.698	0.621
2-Butanone	0.074	0.077	0.070
2,2-Dichloropropane	0.931	0.929	0.854
Cis-1,2-Dichloroethene	0.436	0.435	0.401
Chloroform	0.855	0.847	0.762
Bromochloromethane	0.200	0.194	0.182
1,1,1-Trichloroethane	0.993	0.975	0.872
1,1-Dichloropropene	0.468	0.452	0.424
Carbon Tetrachloride	0.718	0.701	0.637
1,2-Dichloroethane	0.540	0.532	0.492
Benzene	1.178	1.121	1.043
Trichloroethene	0.371	0.360	0.331
1,2-Dichloropropane	0.229	0.218	0.202
Bromodichloromethane	0.497	0.481	0.447
Dibromomethane	0.176	0.169	0.158
2-Chloroethyl Vinyl Ether	0.115	0.113	0.109
4-Methyl-2-Pentanone	0.057	0.054	0.050
Cis 1,3-dichloropropene	0.503	0.490	0.458
Toluene	0.830	0.800	0.734
Trans 1,3-Dichloropropene	0.525	0.499	0.455
2-Hexanone	0.102	0.080	

FORM VI VOA

TN19 : 000055

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

LAB FILE ID: RF20: 2000727

RF40: 4000727

RF60: 6000727

COMPOUND	RF20	RF40	RF60
1,1,2-Trichloroethane	0.221	0.218	0.204
1,3-Dichloropropane	0.402	0.396	0.376
Tetrachloroethene	0.430	0.409	0.384
Chlorodibromomethane	0.378	0.371	0.347
1,2-Dibromoethane	0.244	0.230	0.219
Chlorobenzene	1.021	0.986	0.894
Ethyl Benzene	0.575	0.540	0.497
1,1,1,2-Tetrachloroethane	0.426	0.411	0.376
m,p-xylene	0.690	0.647	0.578
o-Xylene	0.676	0.648	0.591
Styrene	1.125	1.045	0.944
Bromoform	0.358	0.339	
1,1,2,2-Tetrachloroethane	0.405		
1,2,3-Trichloropropane	0.174		
Trans-1,4-Dichloro 2-Butene	0.184	0.132	
N-Propyl Benzene	3.004	2.929	2.640
Bromobenzene	0.746	0.727	0.648
Isopropyl Benzene	2.818	2.721	2.425
2-Chloro Toluene	2.102	2.003	1.874
4-Chloro Toluene	2.340	2.147	2.131
T-Butyl Benzene	2.101	2.066	1.893
1,3,5-Trimethyl Benzene	2.529	2.467	2.227
1,2,4-Trimethylbenzene	2.608	2.552	2.326
S-Butyl Benzene	2.733	2.654	2.408
4-Isopropyl Toluene	2.494	2.465	2.192
1,3-Dichlorobenzene	1.470		
1,4-Dichlorobenzene	1.506	1.492	1.294
N-Butyl Benzene	1.986	1.974	
1,2-Dichlorobenzene	1.375	1.334	1.156
1,2-Dibromo 3-Chloropropane	0.113	0.105	0.090
1,2,4-Trichlorobenzene	0.766		
Hexachloro 1,3-Butadiene	0.385		
Naphthalene	1.353		
1,2,3-Trichlorobenzene	0.398		
Dichlorodifluoromethane	0.604	0.542	0.459
Methyl tert butyl ether	1.204	1.215	1.114

FORM VI VOA

TN19:00059

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

LAB FILE ID: RF20: 2000727 RF40: 4000727 RF60: 6000727

COMPOUND	RF20	RF40	RF60
d4-1,2-Dichloroethane	0.570	0.588	0.563
d8-Toluene	1.125	1.115	1.108
4-Bromofluorobenzene	0.620	0.597	0.505
d4-1,2-Dichlorobenzene	0.894	0.880	0.839
Dibromofluoromethane	0.415	0.422	0.409

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R^2
Chloromethane	LINR		0.9926
Vinyl Chloride	AVRG	0.384	10.4
Bromomethane	AVRG	0.246	10.5
Chloroethane	AVRG	0.237	12.5
Trichlorofluoromethane	AVRG	0.972	17.8
Acrolein	AVRG	0.033	13.6
112Trichloro122Trifluoroetha	AVRG	0.444	12.3
Acetone	AVRG	0.056	9.7
1,1-Dichloroethene	AVRG	0.377	14.1
Bromoethane	AVRG	0.240	13.4
Iodomethane	AVRG	0.573	18.0
Methylene Chloride	AVRG	0.379	17.6
Acrylonitrile	AVRG	0.066	18.2
Carbon Disulfide	AVRG	1.049	9.7
Trans-1,2-Dichloroethene	AVRG	0.452	8.1
Vinyl Acetate	AVRG	0.451	5.1
1,1-Dichloroethane	AVRG	0.714	7.7
2-Butanone	AVRG	0.081	11.3
2,2-Dichloropropane	AVRG	0.929	6.7
Cis-1,2-Dichloroethene	AVRG	0.450	7.3
Chloroform	AVRG	0.909	11.4
Bromochloromethane	AVRG	0.209	10.5
1,1,1-Trichloroethane	AVRG	1.020	9.6
1,1-Dichloropropene	AVRG	0.474	7.4
Carbon Tetrachloride	AVRG	0.722	8.9
1,2-Dichloroethane	AVRG	0.556	8.6
Benzene	AVRG	1.197	8.0
Trichloroethene	AVRG	0.373	7.0
1,2-Dichloropropane	AVRG	0.233	13.8
Bromodichloromethane	AVRG	0.524	9.6
Dibromomethane	AVRG	0.183	9.9
2-Chloroethyl Vinyl Ether	AVRG	0.112	8.4
4-Methyl-2-Pentanone	AVRG	0.058	8.9
Cis 1,3-dichloropropene	AVRG	0.516	8.8
Toluene	AVRG	0.851	8.1
Trans 1,3-Dichloropropene	AVRG	0.524	8.7
2-Hexanone	AVRG	0.105	14.6

<- Indicates value outside QC limits:
(%RSD < 20% or R^2 > 0.990)

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R ²
1,1,2-Trichloroethane	AVRG	0.234	10.0
1,3-Dichloropropane	AVRG	0.438	11.0
Tetrachloroethene	AVRG	0.450	10.1
Chlorodibromomethane	AVRG	0.393	10.1
1,2-Dibromoethane	AVRG	0.252	9.2
Chlorobenzene	AVRG	1.078	9.5
Ethyl Benzene	AVRG	0.578	8.2
1,1,1,2-Tetrachloroethane	AVRG	0.446	9.3
m,p-xylene	AVRG	0.709	10.6
o-Xylene	AVRG	0.662	9.6
Styrene	AVRG	1.101	10.2
Bromoform	AVRG	0.420	15.3
1,1,2,2-Tetrachloroethane	LINR		0.9985
1,2,3-Trichloropropane	AVRG	0.221	19.2
Trans-1,4-Dichloro 2-Butene	2ORDR		0.9956
N-Propyl Benzene	AVRG	3.242	12.2
Bromobenzene	AVRG	0.845	15.6
Isopropyl Benzene	AVRG	2.927	11.6
2-Chloro Toluene	AVRG	2.262	11.7
4-Chloro Toluene	AVRG	2.509	12.6
T-Butyl Benzene	AVRG	2.201	10.2
1,3,5-Trimethyl Benzene	AVRG	2.655	11.3
1,2,4-Trimethylbenzene	AVRG	2.716	9.8
S-Butyl Benzene	AVRG	2.860	9.9
4-Isopropyl Toluene	AVRG	2.525	9.3
1,3-Dichlorobenzene	AVRG	1.656	11.6
1,4-Dichlorobenzene	AVRG	1.694	15.7
N-Butyl Benzene	AVRG	2.061	7.0
1,2-Dichlorobenzene	AVRG	1.489	13.9
1,2-Dibromo 3-Chloropropane	2ORDR		0.9984
1,2,4-Trichlorobenzene	AVRG	0.883	9.8
Hexachloro 1,3-Butadiene	AVRG	0.485	16.1
Naphthalene	AVRG	1.461	7.1
1,2,3-Trichlorobenzene	2ORDR		0.9972
Dichlorodifluoromethane	AVRG	0.569	10.7
Methyl tert butyl ether	AVRG	1.260	7.3

<- Indicates value outside QC limits:
(%RSD < 20% or R² > 0.990)

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Calibration Date: 07/27/11

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R ²
d4-1,2-Dichloroethane	AVRG	0.577	3.4
d8-Toluene	AVRG	1.118	1.0
4-Bromofluorobenzene	AVRG	0.577	6.0
d4-1,2-Dichlorobenzene	AVRG	0.877	2.2
Dibromofluoromethane	AVRG	0.415	2.5

<- Indicates value outside QC limits:
(%RSD < 20% or R² > 0.990)

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

LAB FILE ID: RF0.2: 00_20921 RF0.5: 00_50921 RF1: 01_00921
 RF2: 02_00921 RF10: 10_00921

COMPOUND	RF0.2	RF0.5	RF1	RF2	RF10
Chloromethane		0.942	0.850	0.880	0.733
Vinyl Chloride	0.625	0.610	0.625	0.659	0.642
Bromomethane			0.323	0.325	0.307
Chloroethane	0.427	0.407	0.388	0.412	0.385
Trichlorofluoromethane	0.798	0.711	0.725	0.783	0.724
Acrolein	0.058	0.054	0.057	0.057	0.061
112Trichloro122Trifluoroetha	0.529	0.449	0.475	0.479	0.440
Acetone		0.231	0.193	0.173	0.148
1,1-Dichloroethene	0.625	0.508	0.473	0.498	0.444
Bromoethane	0.340	0.311	0.329	0.338	0.314
Iodomethane			0.526	0.521	0.551
Methylene Chloride		0.598	0.546	0.550	0.507
Acrylonitrile		0.155	0.158	0.169	0.160
Carbon Disulfide	1.638	1.529	1.519	1.651	1.514
Trans-1,2-Dichloroethene	0.564	0.536	0.523	0.543	0.502
Vinyl Acetate	0.978	0.950	0.997	1.058	1.040
1,1-Dichloroethane	1.007	0.928	0.954	0.988	0.910
2-Butanone	0.252	0.236	0.247	0.257	0.244
2,2-Dichloropropane	0.928	0.805	0.770	0.815	0.778
Cis-1,2-Dichloroethene	0.584	0.561	0.559	0.573	0.534
Chloroform	0.956	0.887	0.938	0.963	0.909
Bromochloromethane	0.265	0.250	0.254	0.272	0.250
1,1,1-Trichloroethane	0.897	0.862	0.869	0.900	0.842
1,1-Dichloropropene	0.509	0.504	0.492	0.522	0.490
Carbon Tetrachloride	0.456	0.431	0.439	0.462	0.459
1,2-Dichloroethane	0.498	0.505	0.489	0.510	0.455
Benzene	1.487	1.436	1.433	1.489	1.356
Trichloroethene	0.407	0.350	0.365	0.370	0.346
1,2-Dichloropropane	0.383	0.348	0.350	0.373	0.339
Bromodichloromethane	0.450	0.418	0.421	0.454	0.447
Dibromomethane	0.205	0.183	0.200	0.207	0.187
2-Chloroethyl Vinyl Ether		0.117	0.114	0.107	0.146
4-Methyl-2-Pentanone	0.118	0.126	0.131	0.140	0.135
Cis 1,3-dichloropropene	0.547	0.501	0.535	0.563	0.552
Toluene	0.946	0.932	0.930	0.961	0.874
Trans 1,3-Dichloropropene	0.468	0.467	0.480	0.522	0.521
2-Hexanone	0.251	0.252	0.262	0.286	0.280

FORM VI VOA

TN42:000061

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

LAB FILE ID: RF0.2: 00_20921 RF0.5: 00_50921 RF1: 01_00921
 RF2: 02_00921 RF10: 10_00921

COMPOUND	RF0.2	RF0.5	RF1	RF2	RF10
1,1,2-Trichloroethane	0.328	0.290	0.297	0.308	0.286
1,3-Dichloropropane	0.575	0.562	0.574	0.590	0.558
Tetrachloroethene	0.412	0.403	0.418	0.415	0.394
Chlorodibromomethane	0.263	0.289	0.292	0.317	0.329
1,2-Dibromoethane	0.286	0.288	0.286	0.320	0.301
Chlorobenzene	1.135	1.029	1.070	1.096	1.005
Ethyl Benzene	0.608	0.561	0.552	0.588	0.548
1,1,1,2-Tetrachloroethane	0.347	0.304	0.329	0.358	0.353
m,p-xylene	0.689	0.688	0.693	0.737	0.677
o-Xylene	0.662	0.663	0.679	0.716	0.659
Styrene	0.991	1.031	1.056	1.187	1.121
Bromoform	0.280	0.276	0.289	0.304	0.343
1,1,2,2-Tetrachloroethane	0.602	0.660	0.667	0.683	0.684
1,2,3-Trichloropropane	0.250	0.226	0.232	0.229	0.226
Trans-1,4-Dichloro 2-Butene			0.119	0.136	0.125
N-Propyl Benzene	3.640	3.514	3.562	3.710	3.528
Bromobenzene	0.829	0.749	0.778	0.784	0.746
Isopropyl Benzene	2.501	2.530	2.531	2.670	2.579
2-Chloro Toluene	2.555	2.513	2.509	2.536	2.393
4-Chloro Toluene	2.419	2.278	2.373	2.345	2.236
T-Butyl Benzene	2.417	2.330	2.302	2.437	2.316
1,3,5-Trimethyl Benzene	2.501	2.530	2.531	2.670	2.579
1,2,4-Trimethylbenzene	2.527	2.071	2.293	2.468	2.448
S-Butyl Benzene	3.444	3.038	3.114	3.397	3.276
4-Isopropyl Toluene	2.861	2.170	2.452	2.708	2.665
1,3-Dichlorobenzene	1.651	1.547	1.536	1.610	1.506
1,4-Dichlorobenzene	1.674	1.654	1.641	1.681	1.556
N-Butyl Benzene	2.395	1.518	1.725	2.134	2.087
1,2-Dichlorobenzene	1.615	1.556	1.570	1.588	1.471
1,2-Dibromo 3-Chloropropane		0.112	0.127	0.133	0.135
1,2,4-Trichlorobenzene	1.082	0.757	0.815	0.997	0.996
Hexachloro 1,3-Butadiene			0.559	0.560	0.600
Naphthalene			1.645	1.739	2.138
1,2,3-Trichlorobenzene			0.704	0.719	0.921
Dichlorodifluoromethane	0.428	0.483	0.425	0.502	0.483
Methyl tert butyl ether	1.458	1.374	1.459	1.378	1.400

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

LAB FILE ID: RF0.2: 00_20921 RF0.5: 00_50921 RF1: 01_00921
 RF2: 02_00921 RF10: 10_00921

COMPOUND	RF0.2	RF0.5	RF1	RF2	RF10
d4-1,2-Dichloroethane	0.582	0.575	0.582	0.580	0.575
d8-Toluene	1.240	1.232	1.232	1.241	1.212
4-Bromofluorobenzene	0.558	0.551	0.550	0.557	0.548
d4-1,2-Dichlorobenzene	0.906	0.898	0.914	0.904	0.910
Dibromofluoromethane	0.449	0.444	0.447	0.453	0.459

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

LAB FILE ID: RF20: 20_00921 RF40: 40_00921 RF60: 60_00921

COMPOUND	RF20	RF40	RF60
Chloromethane	0.729	0.758	0.786
Vinyl Chloride	0.640	0.684	0.645
Bromomethane	0.311	0.341	0.355
Chloroethane	0.329		
Trichlorofluoromethane	0.725	0.682	0.628
Acrolein	0.070	0.071	0.073
112Trichloro122Trifluoroetha	0.460	0.467	0.457
Acetone	0.156	0.155	0.157
1,1-Dichloroethene	0.457	0.462	0.460
Bromoethane	0.316	0.313	0.311
Iodomethane	0.545	0.517	0.534
Methylene Chloride	0.524	0.536	0.534
Acrylonitrile	0.172	0.176	0.178
Carbon Disulfide	1.611	1.623	1.614
Trans-1,2-Dichloroethene	0.526	0.528	0.532
Vinyl Acetate	1.115	1.160	1.165
1,1-Dichloroethane	0.954	0.972	0.980
2-Butanone	0.265	0.270	
2,2-Dichloropropane	0.814	0.826	0.822
Cis-1,2-Dichloroethene	0.558	0.567	0.565
Chloroform	0.952	0.968	0.972
Bromochloromethane	0.258	0.259	0.261
1,1,1-Trichloroethane	0.890	0.900	0.893
1,1-Dichloropropene	0.517	0.519	0.517
Carbon Tetrachloride	0.492	0.499	0.506
1,2-Dichloroethane	0.486	0.485	0.492
Benzene	1.419	1.402	1.387
Trichloroethene	0.362	0.363	0.366
1,2-Dichloropropane	0.358	0.361	0.364
Bromodichloromethane	0.488	0.492	0.500
Dibromomethane	0.199	0.198	0.199
2-Chloroethyl Vinyl Ether	0.158	0.162	0.173
4-Methyl-2-Pentanone	0.147	0.144	0.147
Cis 1,3-dichloropropene	0.604	0.610	0.616
Toluene	0.914	0.897	0.892
Trans 1,3-Dichloropropene	0.566	0.576	0.582
2-Hexanone	0.296	0.290	0.279

FORM VI VOA

TN19 : 000064

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

LAB FILE ID: RF20: 20_00921 RF40: 40_00921 RF60: 60_00921

COMPOUND	RF20	RF40	RF60
1,1,2-Trichloroethane	0.304	0.301	0.304
1,3-Dichloropropane	0.575	0.587	0.583
Tetrachloroethene	0.409	0.410	0.404
Chlorodibromomethane	0.360	0.377	0.379
1,2-Dibromoethane	0.318	0.320	0.326
Chlorobenzene	1.037	1.035	1.023
Ethyl Benzene	0.569	0.573	0.571
1,1,1,2-Tetrachloroethane	0.370	0.380	0.382
m,p-xylene	0.700	0.691	0.678
o-Xylene	0.682	0.690	0.695
Styrene	1.186	1.225	1.221
Bromoform	0.409	0.422	0.429
1,1,2,2-Tetrachloroethane	0.728	0.727	0.728
1,2,3-Trichloropropane	0.239	0.234	0.231
Trans-1,4-Dichloro 2-Butene	0.178	0.194	0.200
N-Propyl Benzene	3.624	3.444	3.194
Bromobenzene	0.780	0.778	0.774
Isopropyl Benzene	2.668	2.616	2.475
2-Chloro Toluene	2.494	2.461	2.388
4-Chloro Toluene	2.320	2.283	2.204
T-Butyl Benzene	2.394	2.368	2.271
1,3,5-Trimethyl Benzene	2.668	2.616	2.475
1,2,4-Trimethylbenzene	2.601	2.622	2.505
S-Butyl Benzene	3.306	3.277	3.045
4-Isopropyl Toluene	2.793	2.859	2.682
1,3-Dichlorobenzene	1.565	1.555	1.525
1,4-Dichlorobenzene	1.619	1.606	1.569
N-Butyl Benzene	2.079	2.322	2.111
1,2-Dichlorobenzene	1.524	1.514	1.485
1,2-Dibromo 3-Chloropropane	0.153	0.155	0.158
1,2,4-Trichlorobenzene	0.955	1.014	0.957
Hexachloro 1,3-Butadiene	0.572	0.560	0.548
Naphthalene	1.962	1.832	1.857
1,2,3-Trichlorobenzene	0.865	0.888	0.848
Dichlorodifluoromethane	0.488	0.513	0.472
Methyl tert butyl ether	1.496	1.501	1.544

FORM VI VOA

TN42 : 000005

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

LAB FILE ID: RF20: 20_00921 RF40: 40_00921 RF60: 60_00921

COMPOUND	RF20	RF40	RF60
d4-1,2-Dichloroethane	0.581	0.588	0.588
d8-Toluene	1.242	1.215	1.221
4-Bromofluorobenzene	0.550	0.558	0.573
d4-1,2-Dichlorobenzene	0.918	0.909	0.903
Dibromofluoromethane	0.461	0.470	0.478

FORM VI VOA

TN429 : 2000000

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R ²
Chloromethane	AVRG	0.811	10.0
Vinyl Chloride	AVRG	0.641	3.5
Bromomethane	AVRG	0.327	5.6
Chloroethane	AVRG	0.391	8.7
Trichlorofluoromethane	AVRG	0.722	7.4
Acrolein	AVRG	0.063	11.9
112Trichloro122Trifluoroetha	AVRG	0.470	5.8
Acetone	AVRG	0.173	17.2
1,1-Dichloroethene	AVRG	0.491	11.9
Bromoethane	AVRG	0.321	3.8
Iodomethane	AVRG	0.532	2.6
Methylene Chloride	AVRG	0.542	5.3
Acrylonitrile	AVRG	0.167	5.4
Carbon Disulfide	AVRG	1.587	3.6
Trans-1,2-Dichloroethene	AVRG	0.532	3.3
Vinyl Acetate	AVRG	1.058	7.8
1,1-Dichloroethane	AVRG	0.962	3.3
2-Butanone	AVRG	0.253	4.7
2,2-Dichloropropane	AVRG	0.820	5.9
Cis-1,2-Dichloroethene	AVRG	0.563	2.6
Chloroform	AVRG	0.943	3.2
Bromochloromethane	AVRG	0.259	3.0
1,1,1-Trichloroethane	AVRG	0.882	2.4
1,1-Dichloropropene	AVRG	0.509	2.4
Carbon Tetrachloride	AVRG	0.468	6.0
1,2-Dichloroethane	AVRG	0.490	3.4
Benzene	AVRG	1.426	3.2
Trichloroethene	AVRG	0.366	5.0
1,2-Dichloropropane	AVRG	0.360	3.9
Bromodichloromethane	AVRG	0.459	6.9
Dibromomethane	AVRG	0.197	4.1
2-Chloroethyl Vinyl Ether	AVRG	0.140	19.1
4-Methyl-2-Pentanone	AVRG	0.136	7.7
Cis 1,3-dichloropropene	AVRG	0.566	7.2
Toluene	AVRG	0.918	3.2
Trans 1,3-Dichloropropene	AVRG	0.523	9.2
2-Hexanone	AVRG	0.274	6.4

<- Indicates value outside QC limits:
(%RSD < 20% or R² > 0.990)

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R ²
1,1,2-Trichloroethane	AVRG	0.302	4.2
1,3-Dichloropropane	AVRG	0.576	1.9
Tetrachloroethene	AVRG	0.408	1.8
Chlorodibromomethane	AVRG	0.326	13.3
1,2-Dibromoethane	AVRG	0.306	5.6
Chlorobenzene	AVRG	1.054	4.1
Ethyl Benzene	AVRG	0.571	3.4
1,1,1,2-Tetrachloroethane	AVRG	0.353	7.4
m,p-xylene	AVRG	0.694	2.7
o-Xylene	AVRG	0.681	2.9
Styrene	AVRG	1.127	8.1
Bromoform	AVRG	0.344	19.3
1,1,2,2-Tetrachloroethane	AVRG	0.685	6.4
1,2,3-Trichloropropane	AVRG	0.233	3.4
Trans-1,4-Dichloro 2-Butene	2ORDR		0.9976
N-Propyl Benzene	AVRG	3.527	4.5
Bromobenzene	AVRG	0.777	3.3
Isopropyl Benzene	AVRG	2.571	2.9
2-Chloro Toluene	AVRG	2.481	2.5
4-Chloro Toluene	AVRG	2.307	3.1
T-Butyl Benzene	AVRG	2.354	2.5
1,3,5-Trimethyl Benzene	AVRG	2.571	2.9
1,2,4-Trimethylbenzene	AVRG	2.442	7.4
S-Butyl Benzene	AVRG	3.237	4.8
4-Isopropyl Toluene	AVRG	2.649	8.8
1,3-Dichlorobenzene	AVRG	1.562	3.0
1,4-Dichlorobenzene	AVRG	1.625	2.8
N-Butyl Benzene	AVRG	2.046	14.2
1,2-Dichlorobenzene	AVRG	1.540	3.3
1,2-Dibromo 3-Chloropropane	AVRG	0.139	12.2
1,2,4-Trichlorobenzene	AVRG	0.947	11.4
Hexachloro 1,3-Butadiene	AVRG	0.568	3.0
Naphthalene	AVRG	1.904	10.1
1,2,3-Trichlorobenzene	AVRG	0.833	10.3
Dichlorodifluoromethane	AVRG	0.474	6.8
Methyl tert butyl ether	AVRG	1.451	4.3

<- Indicates value outside QC limits:
(%RSD < 20% or R² > 0.990)

FORM 6
VOLATILE INITIAL CALIBRATION DATA

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Calibration Date: 09/21/11

COMPOUND	CURVE TYPE	AVE RF	%RSD OR R^2
d4-1,2-Dichloroethane	AVRG	0.581	0.9
d8-Toluene	AVRG	1.229	1.0
4-Bromofluorobenzene	AVRG	0.556	1.5
d4-1,2-Dichlorobenzene	AVRG	0.908	0.7
Dibromofluoromethane	AVRG	0.458	2.6

<- Indicates value outside QC limits:
(%RSD < 20% or R^2 > 0.990)

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Cont. Calib. Date: 09/21/11

Init. Calib. Date: 07/27/11

Cont. Calib. Time: 0855

COMPOUND	Cal Amt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
Chloromethane	10.000	10.543	0.100	LINR	5.4
Vinyl Chloride	0.384	0.3528	0.010	AVRG	-8.1
Bromomethane	0.246	0.2324	0.010	AVRG	-5.5
Chloroethane	0.237	0.2146	0.010	AVRG	-9.4
Trichlorofluoromethane	0.972	0.9913	0.010	AVRG	2.0
Acrolein	0.033	0.0303	0.010	AVRG	-8.2
112Trichloro122Trifluoroetha	0.444	0.4240	0.010	AVRG	-4.5
Acetone	0.056	0.0489	0.010	AVRG	-12.7
1,1-Dichloroethene	0.377	0.3766	0.010	AVRG	-0.1
Bromoethane	0.240	0.2370	0.010	AVRG	-1.2
Iodomethane	0.573	0.4921	0.010	AVRG	-14.1
Methylene Chloride	0.379	0.3484	0.010	AVRG	-8.1
Acrylonitrile	0.066	0.0676	0.010	AVRG	2.4
Carbon Disulfide	1.049	0.9808	0.010	AVRG	-6.5
Trans-1,2-Dichloroethene	0.452	0.4315	0.010	AVRG	-4.5
Vinyl Acetate	0.451	0.4325	0.010	AVRG	-4.1
1,1-Dichloroethane	0.714	0.6726	0.100	AVRG	-5.8
2-Butanone	0.081	0.0755	0.010	AVRG	-6.8
2,2-Dichloropropane	0.929	0.9488	0.010	AVRG	2.1
Cis-1,2-Dichloroethene	0.450	0.4126	0.010	AVRG	-8.3
Chloroform	0.909	0.8849	0.010	AVRG	-2.6
Bromochloromethane	0.209	0.2127	0.010	AVRG	1.8
1,1,1-Trichloroethane	1.020	1.0202	0.010	AVRG	0.0
1,1-Dichloropropene	0.474	0.4458	0.010	AVRG	-5.9
Carbon Tetrachloride	0.722	0.7128	0.010	AVRG	-1.3
1,2-Dichloroethane	0.556	0.5306	0.010	AVRG	-4.6
Benzene	1.197	1.1563	0.010	AVRG	-3.4
Trichloroethene	0.373	0.3559	0.010	AVRG	-4.6
1,2-Dichloropropane	0.233	0.2231	0.010	AVRG	-4.2
Bromodichloromethane	0.524	0.5080	0.010	AVRG	-3.0
Dibromomethane	0.183	0.1727	0.010	AVRG	-5.6
2-Chloroethyl Vinyl Ether	0.112	0.1079	0.010	AVRG	-3.7
4-Methyl-2-Pentanone	0.058	0.0562	0.010	AVRG	-3.1
Cis 1,3-dichloropropene	0.516	0.5221	0.010	AVRG	1.2
Toluene	0.851	0.8520	0.010	AVRG	0.1
Trans 1,3-Dichloropropene	0.524	0.5342	0.010	AVRG	1.9
2-Hexanone	0.106	0.1038	0.010	AVRG	-2.1

<- Exceeds QC limit of 20% D
* RF less than minimum RF

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Cont. Calib. Date: 09/21/11

Init. Calib. Date: 07/27/11

Cont. Calib. Time: 0855

COMPOUND	Cal Amt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
1,1,2-Trichloroethane	0.234	0.2259	0.010	AVRG	-3.5
1,3-Dichloropropane	0.438	0.4065	0.010	AVRG	-7.2
Tetrachloroethene	0.450	0.4315	0.010	AVRG	-4.1
Chlorodibromomethane	0.393	0.3813	0.010	AVRG	-3.0
1,2-Dibromoethane	0.252	0.2334	0.010	AVRG	-7.4
Chlorobenzene	1.078	1.0625	0.300	AVRG	-1.4
Ethyl Benzene	0.578	0.5836	0.010	AVRG	1.0
1,1,1,2-Tetrachloroethane	0.445	0.4498	0.010	AVRG	1.1
m,p-xylene	0.709	0.7255	0.010	AVRG	2.3
o-Xylene	0.662	0.6936	0.010	AVRG	4.8
Styrene	1.101	1.1331	0.010	AVRG	2.9
Bromoform	0.420	0.4161	0.100	AVRG	-0.9
1,1,2,2-Tetrachloroethane	10.000	9.575	0.300	LINR	-4.2
1,2,3-Trichloropropane	0.221	0.1824	0.010	AVRG	-17.5
Trans-1,4-Dichloro 2-Butene	10.000	9.288	0.010	2ORDR	-7.1
N-Propyl Benzene	3.242	3.2580	0.010	AVRG	0.5
Bromobenzene	0.845	0.8105	0.010	AVRG	-4.1
Isopropyl Benzene	2.927	3.0401	0.010	AVRG	3.9
2-Chloro Toluene	2.262	2.2173	0.010	AVRG	-2.0
4-Chloro Toluene	2.509	2.4414	0.010	AVRG	-2.7
T-Butyl Benzene	2.201	2.2160	0.010	AVRG	0.7
1,3,5-Trimethyl Benzene	2.655	2.6918	0.010	AVRG	1.4
1,2,4-Trimethylbenzene	2.716	2.7018	0.010	AVRG	-0.5
S-Butyl Benzene	2.860	2.8260	0.010	AVRG	-1.2
4-Isopropyl Toluene	2.525	2.5807	0.010	AVRG	2.2
1,3-Dichlorobenzene	1.656	1.5027	0.010	AVRG	-9.2
1,4-Dichlorobenzene	1.694	1.5410	0.010	AVRG	-9.0
N-Butyl Benzene	2.061	1.9744	0.010	AVRG	-4.2
1,2-Dichlorobenzene	1.489	1.3406	0.010	AVRG	-10.0
1,2-Dibromo 3-Chloropropane	10.000	8.092	0.010	2ORDR	-19.1
1,2,4-Trichlorobenzene	0.883	0.7558	0.010	AVRG	-14.4
Hexachloro 1,3-Butadiene	0.486	0.3883	0.010	AVRG	-20.1
Naphthalene	1.462	1.2831	0.010	AVRG	-12.2
1,2,3-Trichlorobenzene	10.000	9.332	0.010	2ORDR	-6.7
Dichlorodifluoromethane	0.569	0.4152	0.010	AVRG	-27.0
Methyl tert butyl ether	1.259	1.2289	0.010	AVRG	-2.4

<- Exceeds QC limit of 20% D

* RF less than minimum RF

<-

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT5

Cont. Calib. Date: 09/21/11

Init. Calib. Date: 07/27/11

Cont. Calib. Time: 0855

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
d4-1,2-Dichloroethane	0.577	0.6210	0.010	AVRG	7.6
d8-Toluene	1.118	1.1258	0.010	AVRG	0.7
4-Bromofluorobenzene	0.577	0.5841	0.010	AVRG	1.2
d4-1,2-Dichlorobenzene	0.877	0.8668	0.010	AVRG	-1.2
Dibromofluoromethane	0.415	0.4271	0.010	AVRG	2.9

<- Exceeds QC limit of 20% D

* RF less than minimum RRF

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Cont. Calib. Date: 09/26/11

Init. Calib. Date: 09/21/11

Cont. Calib. Time: 0939

COMPOUND	Cal Amt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
Chloromethane	0.811	0.6624	0.100	AVRG	-18.3
Vinyl Chloride	0.641	0.6043	0.010	AVRG	-5.7
Bromomethane	0.327	0.2645	0.010	AVRG	-19.1
Chloroethane	0.391	0.3501	0.010	AVRG	-10.5
Trichlorofluoromethane	0.722	0.7138	0.010	AVRG	-1.1
Acrolein	0.063	0.0569	0.010	AVRG	-9.7
112Trichloro122Trifluoroetha	0.470	0.4519	0.010	AVRG	-3.8
Acetone	0.173	0.1353	0.010	AVRG	-21.8
1,1-Dichloroethene	0.491	0.4429	0.010	AVRG	-9.8
Bromoethane	0.322	0.2964	0.010	AVRG	-8.0
Iodomethane	0.532	0.3701	0.010	AVRG	-30.4
Methylene Chloride	0.542	0.4977	0.010	AVRG	-8.2
Acrylonitrile	0.167	0.1518	0.010	AVRG	-9.1
Carbon Disulfide	1.587	1.5028	0.010	AVRG	-5.3
Trans-1,2-Dichloroethene	0.532	0.5005	0.010	AVRG	-5.9
Vinyl Acetate	1.058	0.9871	0.010	AVRG	-6.7
1,1-Dichloroethane	0.962	0.8756	0.100	AVRG	-9.0
2-Butanone	0.253	0.2338	0.010	AVRG	-7.6
2,2-Dichloropropane	0.820	0.7671	0.010	AVRG	-6.4
Cis-1,2-Dichloroethene	0.563	0.5241	0.010	AVRG	-6.9
Chloroform	0.943	0.8900	0.010	AVRG	-5.6
Bromochloromethane	0.259	0.2441	0.010	AVRG	-5.8
1,1,1-Trichloroethane	0.882	0.8295	0.010	AVRG	-6.0
1,1-Dichloropropene	0.509	0.4885	0.010	AVRG	-4.0
Carbon Tetrachloride	0.468	0.4581	0.010	AVRG	-2.1
1,2-Dichloroethane	0.490	0.4422	0.010	AVRG	-9.8
Benzene	1.426	1.3452	0.010	AVRG	-5.7
Trichloroethene	0.366	0.3516	0.010	AVRG	-3.9
1,2-Dichloropropane	0.360	0.3312	0.010	AVRG	-8.0
Bromodichloromethane	0.459	0.4406	0.010	AVRG	-4.0
Dibromomethane	0.197	0.1840	0.010	AVRG	-6.6
2-Chloroethyl Vinyl Ether	0.140	0.1201	0.010	AVRG	-14.2
4-Methyl-2-Pentanone	0.136	0.1282	0.010	AVRG	-5.7
Cis 1,3-dichloropropene	0.566	0.5365	0.010	AVRG	-5.2
Toluene	0.918	0.8681	0.010	AVRG	-5.4
Trans 1,3-Dichloropropene	0.523	0.5003	0.010	AVRG	-4.3
2-Hexanone	0.274	0.2542	0.010	AVRG	-7.2

<- Exceeds QC limit of 20% D

* RF less than minimum RF

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Cont. Calib. Date: 09/26/11

Init. Calib. Date: 09/21/11

Cont. Calib. Time: 0939

COMPOUND	Cal Amt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
1,1,2-Trichloroethane	0.302	0.2814	0.010	AVRG	-6.8
1,3-Dichloropropane	0.576	0.5226	0.010	AVRG	-9.3
Tetrachloroethene	0.408	0.4040	0.010	AVRG	-1.0
Chlorodibromomethane	0.326	0.3289	0.010	AVRG	0.9
1,2-Dibromoethane	0.306	0.2876	0.010	AVRG	-6.0
Chlorobenzene	1.054	0.9954	0.300	AVRG	-5.6
Ethyl Benzene	0.571	0.5373	0.010	AVRG	-5.9
1,1,1,2-Tetrachloroethane	0.353	0.3456	0.010	AVRG	-2.1
m,p-xylene	0.694	0.6739	0.010	AVRG	-2.9
o-Xylene	0.681	0.6435	0.010	AVRG	-5.5
Styrene	1.127	1.1296	0.010	AVRG	0.2
Bromoform	0.344	0.3439	0.100	AVRG	-0.0
1,1,2,2-Tetrachloroethane	0.685	0.6328	0.300	AVRG	-7.6
1,2,3-Trichloropropane	0.233	0.2092	0.010	AVRG	-10.2
Trans-1,4-Dichloro 2-Butene	10.000	7.018	0.010	2ORDR	-29.8
N-Propyl Benzene	3.527	3.3874	0.010	AVRG	-4.0
Bromobenzene	0.777	0.7278	0.010	AVRG	-6.3
Isopropyl Benzene	2.571	2.4754	0.010	AVRG	-3.7
2-Chloro Toluene	2.481	2.2921	0.010	AVRG	-7.6
4-Chloro Toluene	2.307	2.1471	0.010	AVRG	-6.9
T-Butyl Benzene	2.354	2.2400	0.010	AVRG	-4.8
1,3,5-Trimethyl Benzene	2.571	2.4754	0.010	AVRG	-3.7
1,2,4-Trimethylbenzene	2.442	2.4454	0.010	AVRG	0.1
S-Butyl Benzene	3.237	3.1945	0.010	AVRG	-1.3
4-Isopropyl Toluene	2.649	2.7596	0.010	AVRG	4.2
1,3-Dichlorobenzene	1.562	1.4829	0.010	AVRG	-5.1
1,4-Dichlorobenzene	1.625	1.5263	0.010	AVRG	-6.1
N-Butyl Benzene	2.046	2.1562	0.010	AVRG	5.4
1,2-Dichlorobenzene	1.540	1.4386	0.010	AVRG	-6.6
1,2-Dibromo 3-Chloropropane	0.139	0.1218	0.010	AVRG	-12.4
1,2,4-Trichlorobenzene	0.947	0.9954	0.010	AVRG	5.1
Hexachloro 1,3-Butadiene	0.568	0.5674	0.010	AVRG	-0.1
Naphthalene	1.904	2.0082	0.010	AVRG	5.5
1,2,3-Trichlorobenzene	0.833	0.8534	0.010	AVRG	2.4
Dichlorodifluoromethane	0.474	0.4563	0.010	AVRG	-3.7
Methyl tert butyl ether	1.451	1.3215	0.010	AVRG	-8.9

<- Exceeds QC limit of 20% D

* RF less than minimum RF

7A
VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIG

Instrument ID: NT2

Cont. Calib. Date: 09/26/11

Init. Calib. Date: 09/21/11

Cont. Calib. Time: 0939

COMPOUND	CalAmt or ARF	CC Amt or RF	MIN RRF	CURVE TYPE	%D or Drift
d4-1,2-Dichloroethane	0.581	0.5513	0.010	AVRG	-5.1
d8-Toluene	1.229	1.2170	0.010	AVRG	-1.0
4-Bromofluorobenzene	0.556	0.5536	0.010	AVRG	-0.4
d4-1,2-Dichlorobenzene	0.908	0.9103	0.010	AVRG	0.2
Dibromofluoromethane	0.458	0.4536	0.010	AVRG	-1.0

<- Exceeds QC limit of 20% D

* RF less than minimum RF

8A
VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIGA

Ical Midpoint ID: 1000727

Ical Date: 07/27/11

Instrument ID: NT5

Project Run Date: 09/21/11

	IS1 (PFB) AREA #	RT #	IS2 (DFB) AREA #	RT #	IS3 (CLB) AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	319331	4.77	421917	5.21	402459	7.67
UPPER LIMIT	638662	5.27	843834	5.71	804918	8.17
LOWER LIMIT	159666	4.27	210958	4.71	201230	7.17
=====	=====	=====	=====	=====	=====	=====
Sample ID						
=====	=====	=====	=====	=====	=====	=====
01 LCS0921	333608	4.77	454880	5.21	420554	7.67
02 LCS0921	316220	4.77	427131	5.21	395480	7.67
03 MB0921	297550	4.77	419779	5.21	387163	7.66
04 FRP-091911-0	243958	4.77	318525	5.21	289593	7.66
05 FRP-091911-0	259731	4.77	365931	5.21	382933	7.66
06 FRP-091911-0	292481	4.85	462562	5.28	394770	7.75
07 FRP-091911-0	293301	4.77	385524	5.21	364245	7.66
08 FRP-091911-0	265354	4.77	356334	5.21	338365	7.66
09 TRIP BLANKS	258822	4.77	347379	5.21	328343	7.66
10 FRP-092011-0	264434	4.77	356649	5.21	320607	7.67
11 FRP-092011-0	258738	4.77	345053	5.21	317057	7.67
12 TRIP BLANKS	255052	4.77	337989	5.21	315968	7.66
13						
14						
15						
16						
17						
18						
19						
20						
21						
22						

IS1 (PFB) = Pentafluorobenzene

IS2 (DFB) = 1,4-Difluorobenzene

IS3 (CLB) = d5-Chlorobenzene

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint

AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint

RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Ical midpoint

RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Ical midpoint

* Values outside of QC limits.

8A
VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIGA

Ical Midpoint ID: 1000727

Ical Date: 07/27/11

Instrument ID: NT5

Project Run Date: 09/21/11

	IS4 (DCB) AREA #	RT #	AREA #	RT #	AREA #	RT #
ICAL MIDPT	245371	9.72				
UPPER LIMIT	490742	10.22				
LOWER LIMIT	122686	9.22				
Sample ID						
01 LCS0921	251539	9.72				
02 LCS0921	223163	9.72				
03 MB0921	200860	9.72				
04 FRP-091911-0	152885	9.72				
05 FRP-091911-0	206198	9.72				
06 FRP-091911-0	152847	9.76				
07 FRP-091911-0	214077	9.72				
08 FRP-091911-0	186056	9.72				
09 TRIP BLANKS	174317	9.72				
10 FRP-092011-0	179204	9.72				
11 FRP-092011-0	173682	9.72				
12 TRIP BLANKS	167717	9.72				
13						
14						
15						
16						
17						
18						
19						
20						
21						
22						

IS4 (DCB) = d4-1,4-Dichlorobenzene

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint
 AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint
 RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Ical midpoint
 RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Ical midpoint

* Values outside of QC limits.

8A
VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIGA

Ical Midpoint ID: 10_00921

Ical Date: 09/21/11

Instrument ID: NT2

Project Run Date: 09/26/11

	IS1 (PFB) AREA #	RT #	IS2 (DFB) AREA #	RT #	IS3 (CLB) AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	331438	5.46	505777	5.86	491374	7.92
UPPER LIMIT	662876	5.96	1011554	6.36	982748	8.42
LOWER LIMIT	165719	4.96	252888	5.36	245687	7.42
=====	=====	=====	=====	=====	=====	=====
Sample ID						
=====	=====	=====	=====	=====	=====	=====
01 LCS0926	302141	5.46	463384	5.86	458683	7.92
02 LCS0926	294107	5.46	447017	5.86	452687	7.92
03 MB0926	310738	5.46	468089	5.86	469380	7.92
04 TRIP BLANKS	306553	5.46	459320	5.86	459420	7.92
05 FRP-091911-0	321519	5.46	480140	5.86	472582	7.92
06 FRP-091911-0	312530	5.46	469646	5.85	462564	7.92
07 FRP-091911-0	318130	5.46	478761	5.85	472117	7.92
08 FRP-091911-0	312866	5.46	470323	5.85	465816	7.92
09 FRP-092011-0	305691	5.46	455665	5.86	454285	7.92
10 FRP-092011-0	297479	5.46	445846	5.86	448317	7.92
11 FRP-092011-0	304173	5.46	455255	5.86	455236	7.92
12 FRP-092111-0	317417	5.46	475118	5.85	466598	7.92
13 FRP-092111-0	321297	5.46	479180	5.85	472408	7.92
14 FRP-092111-0	318994	5.46	480834	5.85	469531	7.92
15 FRP-092111-0	324022	5.46	490358	5.85	479960	7.92
16 FRP-092111-0	319766	5.46	483678	5.85	474939	7.92
17 FRP-092111-0	299278	5.46	446469	5.86	445768	7.92
18 FRP-091911-0	315458	5.46	478667	5.85	464572	7.92
19 FRP-091911-0	324929	5.46	492284	5.85	477416	7.92
20						
21						
22						

IS1 (PFB) = Pentafluorobenzene

IS2 (DFB) = 1,4-Difluorobenzene

IS3 (CLB) = d5-Chlorobenzene

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint

AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint

RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Ical midpoint

RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Ical midpoint

* Values outside of QC limits.

8A
VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: ANALYTICAL RESOURCES INC

Client: AMEC EARTH AND ENVIRONMENTAL

ARI Job No: TN42

Project: 2011 FRP SHORELINE INVESTIGA

Ical Midpoint ID: 10_00921

Ical Date: 09/21/11

Instrument ID: NT2

Project Run Date: 09/26/11

	IS4 (DCB) AREA #	RT #	AREA #	RT #	AREA #	RT #
=====	=====	=====	=====	=====	=====	=====
ICAL MIDPT	306350	9.62				
UPPER LIMIT	612700	10.12				
LOWER LIMIT	153175	9.12				
=====	=====	=====	=====	=====	=====	=====
Sample ID						
=====	=====	=====	=====	=====	=====	=====
01 LCS0926	286825	9.61				
02 LCS0926	286948	9.61				
03 MB0926	284610	9.61				
04 TRIP BLANKS	282436	9.61				
05 FRP-091911-0	286008	9.61				
06 FRP-091911-0	277671	9.61				
07 FRP-091911-0	278594	9.61				
08 FRP-091911-0	277588	9.61				
09 FRP-092011-0	266535	9.61				
10 FRP-092011-0	273338	9.61				
11 FRP-092011-0	277228	9.61				
12 FRP-092111-0	278502	9.61				
13 FRP-092111-0	275481	9.61				
14 FRP-092111-0	270111	9.61				
15 FRP-092111-0	277630	9.61				
16 FRP-092111-0	279134	9.61				
17 FRP-092111-0	275086	9.61				
18 FRP-091911-0	284602	9.61				
19 FRP-091911-0	279507	9.61				
20						
21						
22						

IS4 (DCB) = d4-1,4-Dichlorobenzene

AREA UPPER LIMIT = +100% of internal standard area from Ical midpoint

AREA LOWER LIMIT = - 50% of internal standard area from Ical midpoint

RT UPPER LIMIT = + 0.50 minutes of internal standard RT from Ical midpoint

RT LOWER LIMIT = - 0.50 minutes of internal standard RT from Ical midpoint

* Values outside of QC limits.

Metals Analysis
Report and Summary QC Forms

ARI Job ID: TN19, TN21

Cover Page**INORGANIC ANALYSIS DATA PACKAGE****ANALYTICAL
RESOURCES
INCORPORATED**

CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

SDG: TN19

CLIENT ID	ARI ID	ARI LIMS ID	REPREP
FRP-092011-001	TN19A	11-20545	
FRP-092011-001D	TN19ADUP	11-20545	
FRP-092011-001S	TN19ASPK	11-20545	
FRP-092011-002	TN19B	11-20546	
PBW	TN19MB1	11-20546	
LCSW	TN19MB1SPK	11-20546	
FRP-092011-003	TN19C	11-20547	
FRP-092011-004	TN19D	11-20548	
FRP-092011-005	TN19E	11-20549	

Were ICP interelement corrections applied ? Yes/No YES

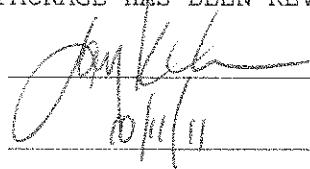
Were ICP background corrections applied ? Yes/No YES

If yes - were raw data generated before application of background corrections ? Yes/No NO

Comments:

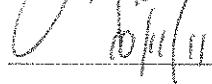
THIS DATA PACKAGE HAS BEEN REVIEWED AND AUTHORIZED FOR RELEASE BY:

Signature:



Name: Jay Kuhn

Date:



Title: Inorganics Director

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: FRP-092011-001
SAMPLE

Lab Sample ID: TN19A
LIMS ID: 11-20545
Matrix: Water
Data Release Authorized:
Reported: 10/11/11

QC Report No: TN19-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769
Date Sampled: 09/20/11
Date Received: 09/20/11

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	MDL	RL	Result	Q
3010A	09/22/11	6010B	09/27/11	7429-90-5	Aluminum	0.128	0.2	880	
200.8	09/22/11	200.8	10/07/11	7440-38-2	Arsenic	0.00024	0.001	0.096	
3010A	09/22/11	6010B	09/27/11	7440-43-9	Cadmium	0.0009	0.01	0.01	U
3010A	09/22/11	6010B	09/27/11	7440-47-3	Chromium	0.0062	0.02	0.85	
3010A	09/22/11	6010B	09/27/11	7440-50-8	Copper	0.0046	0.01	1.36	
200.8	09/22/11	200.8	10/07/11	7439-92-1	Lead	0.000230	0.0005	0.0782	
3010A	09/22/11	6010B	09/27/11	7440-02-0	Nickel	0.0193	0.05	0.51	
3010A	09/22/11	6010B	09/27/11	7782-49-2	Selenium	0.025	0.2	0.2	U
3010A	09/22/11	6010B	09/27/11	7440-28-0	Thallium	0.016	0.2	0.2	U
3010A	09/22/11	6010B	09/27/11	7440-62-2	Vanadium	0.0014	0.02	2.22	
3010A	09/22/11	6010B	09/27/11	7440-66-6	Zinc	0.0072	0.05	1.80	

Reported in mg/L (ppm).

U-Analyte undetected at given RL

RL=Reporting Limit

INORGANICS ANALYSIS DATA SHEET
TOTAL METALS

Page 1 of 1

Sample ID: FRP-092011-002

SAMPLE

Lab Sample ID: TN19B

LIMS ID: 11-20546

Matrix: Water

Data Release Authorized: ✓

Reported: 10/11/11

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation

8769

Date Sampled: 09/20/11

Date Received: 09/20/11

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	MDL	RL	Result	Q
3010A	09/22/11	6010B	09/27/11	7429-90-5	Aluminum	0.0257	0.05	10.7	
200.8	09/22/11	200.8	10/06/11	7440-38-2	Arsenic	0.000120	0.0005	0.0086	
3010A	09/22/11	6010B	09/27/11	7440-43-9	Cadmium	0.00018	0.002	0.002	U
3010A	09/22/11	6010B	09/27/11	7440-47-3	Chromium	0.00124	0.005	0.039	
3010A	09/22/11	6010B	09/27/11	7440-50-8	Copper	0.00092	0.002	0.083	
200.8	09/22/11	200.8	10/06/11	7439-92-1	Lead	0.000115	0.0002	0.0064	
3010A	09/22/11	6010B	09/27/11	7440-02-0	Nickel	0.0039	0.01	0.01	U
3010A	09/22/11	6010B	09/27/11	7782-49-2	Selenium	0.0050	0.05	0.05	U
3010A	09/22/11	6010B	09/27/11	7440-28-0	Thallium	0.0031	0.05	0.05	U
3010A	09/22/11	6010B	09/27/11	7440-62-2	Vanadium	0.00027	0.003	0.200	
3010A	09/22/11	6010B	09/27/11	7440-66-6	Zinc	0.0014	0.01	0.02	

Reported in mg/L (ppm).

U-Analyte undetected at given RL

RL=Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: FRP-092011-003

SAMPLE

Lab Sample ID: TN19C

LIMS ID: 11-20547

Matrix: Water

Data Release Authorized:

Reported: 10/11/11

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation

8769

Date Sampled: 09/20/11

Date Received: 09/20/11

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	MDL	RL	Result	Q
3010A	09/22/11	6010B	09/27/11	7429-90-5	Aluminum	0.0257	0.05	18.9	
200.8	09/22/11	200.8	10/06/11	7440-38-2	Arsenic	0.000120	0.0005	0.0152	
3010A	09/22/11	6010B	09/27/11	7440-43-9	Cadmium	0.00018	0.002	0.002	U
3010A	09/22/11	6010B	09/27/11	7440-47-3	Chromium	0.00124	0.005	0.068	
3010A	09/22/11	6010B	09/27/11	7440-50-8	Copper	0.00092	0.002	0.072	
200.8	09/22/11	200.8	10/06/11	7439-92-1	Lead	0.000115	0.0002	0.0072	
3010A	09/22/11	6010B	09/27/11	7440-02-0	Nickel	0.0039	0.01	0.01	U
3010A	09/22/11	6010B	09/27/11	7782-49-2	Selenium	0.0050	0.05	0.05	U
3010A	09/22/11	6010B	09/27/11	7440-28-0	Thallium	0.0031	0.05	0.05	U
3010A	09/22/11	6010B	09/27/11	7440-62-2	Vanadium	0.00027	0.003	0.315	
3010A	09/22/11	6010B	09/27/11	7440-66-6	Zinc	0.0014	0.01	0.02	

Reported in mg/L (ppm).

U-Analyte undetected at given RL

RL=Reporting Limit

CMW

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: FRP-092011-004

SAMPLE

Lab Sample ID: TN19D

LIMS ID: 11-20548

Matrix: Water

Data Release Authorized:

Reported: 10/11/11

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation

8769

Date Sampled: 09/20/11

Date Received: 09/20/11

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	MDL	RL	Result	Q
3010A	09/22/11	6010B	09/27/11	7429-90-5	Aluminum	0.0257	0.05	82.9	
200.8	09/22/11	200.8	10/06/11	7440-38-2	Arsenic	0.000120	0.0005	0.0249	
3010A	09/22/11	6010B	09/27/11	7440-43-9	Cadmium	0.00018	0.002	0.002	U
3010A	09/22/11	6010B	09/27/11	7440-47-3	Chromium	0.00124	0.005	0.258	
3010A	09/22/11	6010B	09/27/11	7440-50-8	Copper	0.00092	0.002	0.182	
200.8	09/22/11	200.8	10/06/11	7439-92-1	Lead	0.000115	0.0002	0.0128	
3010A	09/22/11	6010B	09/27/11	7440-02-0	Nickel	0.0039	0.01	0.11	
3010A	09/22/11	6010B	09/27/11	7782-49-2	Selenium	0.0050	0.05	0.05	U
3010A	09/22/11	6010B	09/27/11	7440-28-0	Thallium	0.0031	0.05	0.05	U
3010A	09/22/11	6010B	09/27/11	7440-62-2	Vanadium	0.00027	0.003	0.361	
3010A	09/22/11	6010B	09/27/11	7440-66-6	Zinc	0.0014	0.01	0.33	

Reported in mg/L (ppm).

U-Analyte undetected at given RL

RL=Reporting Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: FRP-092011-005

SAMPLE

Lab Sample ID: TN19E

LIMS ID: 11-20549

Matrix: Water

Data Release Authorized:

Reported: 10/11/11

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation

8769

Date Sampled: 09/20/11

Date Received: 09/20/11

Prep Meth	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	MDL	RL	Result	Q
3010A	09/22/11	6010B	09/27/11	7429-90-5	Aluminum	0.0257	0.05	0.05	U
200.8	09/22/11	200.8	10/07/11	7440-38-2	Arsenic	0.000048	0.0002	0.0002	U
3010A	09/22/11	6010B	09/27/11	7440-43-9	Cadmium	0.00018	0.002	0.002	U
3010A	09/22/11	6010B	09/27/11	7440-47-3	Chromium	0.00124	0.005	0.005	U
3010A	09/22/11	6010B	09/27/11	7440-50-8	Copper	0.00092	0.002	0.002	U
200.8	09/22/11	200.8	10/07/11	7439-92-1	Lead	0.000046	0.0001	0.0001	U
3010A	09/22/11	6010B	09/27/11	7440-02-0	Nickel	0.0039	0.01	0.01	U
3010A	09/22/11	6010B	09/27/11	7782-49-2	Selenium	0.0050	0.05	0.05	U
3010A	09/22/11	6010B	09/27/11	7440-28-0	Thallium	0.0031	0.05	0.05	U
3010A	09/22/11	6010B	09/27/11	7440-62-2	Vanadium	0.00027	0.003	0.003	U
3010A	09/22/11	6010B	09/27/11	7440-66-6	Zinc	0.0014	0.01	0.01	U

Reported in mg/L (ppm).

U-Analyte undetected at given RL

RL-Reporting Limit

INORGANICS ANALYSIS DATA SHEET
TOTAL METALS

Page 1 of 1

 Sample ID: FRP-092011-001
MATRIX SPIKE

Lab Sample ID: TN19A

LIMS ID: 11-20545

Matrix: Water

 Data Release Authorized: *[Signature]*

Reported: 10/11/11

QC Report No: TN19-AMEC Geomatrix

 Project: FRP 2011 Shoreline Investigation
 8769

Date Sampled: 09/20/11

Date Received: 09/20/11

MATRIX SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Spike	Spike Added	% Recovery	Q
Aluminum	6010B	880	1,100	2.0	11000%	H
Arsenic	200.8	0.096	0.121	0.025	100%	
Cadmium	6010B	0.010 U	0.409	0.500	81.8%	
Chromium	6010B	0.85	1.35	0.50	100%	
Copper	6010B	1.36	1.86	0.500	100%	
Lead	200.8	0.0782	0.0970	0.0250	75.2%	
Nickel	6010B	0.51	0.94	0.50	86.0%	
Selenium	6010B	0.2 U	1.2	2.0	60.0%	N
Thallium	6010B	0.2 U	1.6	2.0	80.0%	
Vanadium	6010B	2.22	2.82	0.50	120%	H
Zinc	6010B	1.80	2.34	0.50	108%	

Reported in mg/L

N-Control Limit Not Met

H-% Recovery Not Applicable, Sample Concentration Too High

NA-Not Applicable, Analyte Not Spiked

Percent Recovery Limits: 75-125%

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: FRP-092011-001
DUPLICATE

Lab Sample ID: TN19A

LIMS ID: 11-20545

Matrix: Water

Data Release Authorized:

Reported: 10/11/11

QC Report No: TN19-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation
8769

Date Sampled: 09/20/11

Date Received: 09/20/11

MATRIX DUPLICATE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Duplicate	RPD	Control Limit	Q
Aluminum	6010B	880	784	11.5%	+/- 20%	
Arsenic	200.8	0.096	0.111	14.5%	+/- 20%	
Cadmium	6010B	0.01 U	0.01 U	0.0%	+/- 0.01	L
Chromium	6010B	0.85	0.75	12.5%	+/- 20%	
Copper	6010B	1.36	1.13	18.5%	+/- 20%	
Lead	200.8	0.0782	0.0852	8.6%	+/- 20%	
Nickel	6010B	0.51	0.44	14.7%	+/- 20%	
Selenium	6010B	0.2 U	0.2 U	0.0%	+/- 0.2	L
Thallium	6010B	0.2 U	0.2 U	0.0%	+/- 0.2	L
Vanadium	6010B	2.22	1.93	14.0%	+/- 20%	
Zinc	6010B	1.80	1.58	13.0%	+/- 20%	

Reported in mg/L

*-Control Limit Not Met

L-RPD Invalid, Limit = Detection Limit

INORGANICS ANALYSIS DATA SHEET
TOTAL METALS

Page 1 of 1

Sample ID: LAB CONTROL

 Lab Sample ID: TN19LCS
 LIMS ID: 11-20546
 Matrix: Water
 Data Release Authorized: *[Signature]*
 Reported: 10/11/11

 QC Report No: TN19-AMEC Geomatrix
 Project: FRP 2011 Shoreline Investigation
 8769
 Date Sampled: NA
 Date Received: NA

BLANK SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Spike Found	Spike Added	% Recovery	Q
Aluminum	6010B	2.02	2.00	101%	
Arsenic	200.8	0.0259	0.0250	104%	
Cadmium	6010B	0.503	0.500	101%	
Chromium	6010B	0.493	0.500	98.6%	
Copper	6010B	0.495	0.500	99.0%	
Lead	200.8	0.0266	0.0250	106%	
Nickel	6010B	0.47	0.50	94.0%	
Selenium	6010B	1.85	2.00	92.5%	
Thallium	6010B	1.85	2.00	92.5%	
Vanadium	6010B	0.484	0.500	96.8%	
Zinc	6010B	0.47	0.50	94.0%	

Reported in mg/L

 N-Control limit not met
 Control Limits: 80-120%

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: METHOD BLANK

Lab Sample ID: TN19MB
 LIMS ID: 11-20546
 Matrix: Water
 Data Release Authorized:
 Reported: 10/11/11

QC Report No: TN19-AMEC Geomatrix
 Project: FRP 2011 Shoreline Investigation
 8769
 Date Sampled: NA
 Date Received: NA

Prep Method	Prep Date	Analysis Method	Analysis Date	CAS Number	Analyte	MDL	RL	Result	Q
3010A	09/22/11	6010B	09/27/11	7429-90-5	Aluminum	0.0257	0.05	0.05	U
200.8	09/22/11	200.8	10/06/11	7440-38-2	Arsenic	0.000048	0.0002	0.0002	U
3010A	09/22/11	6010B	09/27/11	7440-43-9	Cadmium	0.00018	0.002	0.002	U
3010A	09/22/11	6010B	09/27/11	7440-47-3	Chromium	0.00124	0.005	0.005	U
3010A	09/22/11	6010B	09/27/11	7440-50-8	Copper	0.00092	0.002	0.002	U
200.8	09/22/11	200.8	10/06/11	7439-92-1	Lead	0.000046	0.0001	0.0001	U
3010A	09/22/11	6010B	09/27/11	7440-02-0	Nickel	0.0039	0.01	0.01	U
3010A	09/22/11	6010B	09/27/11	7782-49-2	Selenium	0.0050	0.05	0.05	U
3010A	09/22/11	6010B	09/27/11	7440-28-0	Thallium	0.0031	0.05	0.05	U
3010A	09/22/11	6010B	09/27/11	7440-62-2	Vanadium	0.00027	0.003	0.003	U
3010A	09/22/11	6010B	09/27/11	7440-66-6	Zinc	0.0014	0.01	0.01	U

Reported in mg/L (ppm).

U-Analyte undetected at given RL

RL=Reporting Limit

Calibration Verification

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN19

UNITS: ug/L



ANALYTE	EL	M	RUN	ICVTV	ICV	%R	CCVTv	CCV1	%R	CCV2	%R	CCV3	%R	CCV4	%R	CCV5	%R
Aluminum	AL	ICP	IP092771	2000.0	1988.30	99.4	2000.0	2017.69	100.9	2040.49	102.0	1970.63	98.5	1963.83	98.2	1971.51	98.6
Arsenic	AS	PMS	MS100682	50.0	50.04	100.1	50.0	49.47	98.9	49.11	98.2	49.56	99.1	50.17	100.3	49.55	99.1
Cadmium	CD	ICP	IP092771	1000.0	1009.47	100.9	1000.0	1016.89	101.7	1035.92	103.6	1021.69	102.2	1002.41	100.2	1002.51	100.3
Chromium	CR	ICP	IP092771	1000.0	992.76	99.3	1000.0	997.19	99.7	1017.50	101.8	984.05	98.4	977.84	97.8	984.62	98.5
Copper	CU	ICP	IP092771	1000.0	970.43	97.0	1000.0	980.82	98.1	1000.32	100.0	990.53	99.1	973.04	97.3	966.19	96.6
Lead	PB	PMS	MS100682	50.0	49.54	99.1	50.0	49.07	98.1	48.89	97.8	48.84	97.7	48.99	98.0	49.54	99.1
Nickel	NI	ICP	IP092771	1000.0	987.11	98.7	1000.0	977.68	97.8	992.25	99.2	949.35	94.9	945.56	94.6	950.03	95.0
Selenium	SE	ICP	IP092771	2000.0	1889.99	94.5	2000.0	1882.97	94.1	1895.26	94.8	1850.95	92.5	1809.01	90.5	1803.80	90.2
Thallium	TI	ICP	IP092771	2000.0	1888.64	94.4	2000.0	1880.70	94.0	1897.32	94.9	1840.73	92.0	1800.53	90.0	1795.71	89.8
Vanadium	V	ICP	IP092771	1000.0	936.07	93.6	1000.0	940.86	94.1	959.18	95.9	945.69	94.6	930.59	93.1	927.77	92.8
Zinc	ZN	ICP	IP092771	1000.0	978.61	97.9	1000.0	975.28	97.5	996.03	99.6	962.06	96.2	958.41	95.8	969.58	97.0

Control Limits: Mercury 80-120; Other Metals 90-110

FORM II (1)

Calibration Verification

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN19

UNITS: ug/L



ANALYTE	EL	M	RUN	CCV7V	CCV6 %R	CCV7 %R	CCV8 %R	CCV9 %R	CCV10 %R	CCV11 %R
Aluminum	AL	ICP	IP092771	2000.0	1953.33	97.7	1985.46	99.3	1980.30	99.0
Arsenic	AS	PMS	MS100682	50.0	50.12	100.2	50.83	101.7		
Cadmium	CD	ICP	IP092771	1000.0	1011.85	101.2	1024.67	102.5	1017.58	101.8
Chromium	CR	ICP	IP092771	1000.0	993.29	99.3	1010.51	101.1	1003.72	100.4
Copper	CU	ICP	IP092771	1000.0	964.69	96.5	977.08	97.7	971.83	97.2
Lead	PB	PMS	MS100682	50.0	48.91	97.8	49.44	98.9		
Nickel	NI	ICP	IP092771	1000.0	1012.21	101.2	1032.71	103.3	1028.05	102.8
Selenium	SE	ICP	IP092771	2000.0	2030.04	101.5	2043.80	102.2	2016.20	100.8
Thallium	TL	ICP	IP092771	2000.0	2000.47	100.0	2016.53	100.8	1994.11	99.7
Vanadium	V	ICP	IP092771	1000.0	928.54	92.9	946.09	94.6	939.76	94.0
Zinc	ZN	ICP	IP092771	1000.0	994.84	99.5	1013.46	101.3	1006.94	100.7

Control Limits: Mercury 80-120; Other Metals 90-110

FORM II (1)

Calibration Verification

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN19



UNITS: ug/L

ANALYTE	EL	M	RUN	ICVTV	ICV	%R	CCVTV	CCV1	%R	CCV2	%R	CCV3	%R	CCV4	%R	CCV5	%R
Arsenic	AS	PMS	MS100781	50.0	49.12	98.2	50.0	49.66	99.3	50.34	100.7	50.86	101.7	51.41	102.8	51.67	103.3
Lead	PB	PMS	MS100781	50.0	48.89	97.8	50.0	48.79	97.6	48.49	97.0	47.95	95.9	47.66	95.3	47.86	95.7

100-000000

Control Limits: Mercury 80-120; Other Metals 90-110

FORM II (1)

CRDL Standard

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN19

UNITS: ug/L

ANALYTE	EL	M	RUN	CRA/I	TV	CR-1	%R	CR-2	%R	CR-3	%R	CR-4	%R	CR-5	%R	CR-6	%R
Aluminum	AL	ICP	IP092771	50.0		53.37	106.7	49.58		99.2							
Arsenic	AS	PMS	MS100682	0.2		0.19	95.0										
Cadmium	CD	ICP	IP092771	2.0		2.27	113.5	2.13	106.5								
Chromium	CR	ICP	IP092771	5.0		4.96	99.2	4.57	91.4								
Copper	CU	ICP	IP092771	2.0		2.84	142.0	1.90	95.0								
Lead	PB	PMS	MS100682	0.1		0.11	110.0										
Nickel	NI	ICP	IP092771	10.0		9.00	90.0										
Selenium	SE	ICP	IP092771	50.0		50.50	101.0										
Thallium	TL	ICP	IP092771	50.0		47.14	94.3	48.83	97.7								
Vanadium	V	ICP	IP092771	3.0		3.10	103.3	3.01	100.3								
Zinc	ZN	ICP	IP092771	10.0		9.91	99.1	9.42	94.2								
Arsenic	AS	PMS	MS100781	0.2		0.17	85.0										
Lead	PB	PMS	MS100781	0.1		0.11	110.0										

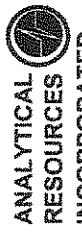
Control Limits: no control limits have been established by the EPA at this time.

Calibration Blanks

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SINGING



UNIT STUDY/T

SDG INDEX

TN19 : 66295

Calibration Blanks

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN19



UNITS: ug/L

ANALYTE	EL	METH	RUN	CRDL	IDL	CCB6	C	CCB7	C	CCB8	C	CCB9	C	CCB10	C	CCB11	C
Aluminum	AL	ICP	IPO92771	200.0	50.0	50.0	U	50.0	U	50.0	U	50.0	U	50.0	U	50.0	U
Arsenic	AS	PMS	MS100682	10.0	0.2	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U
Cadmium	CD	ICP	IPO92771	5.0	2.0	2.0	U	2.0	U	2.0	U	2.0	U	2.0	U	2.0	U
Chromium	CR	ICP	IPO92771	10.0	5.0	5.0	U	5.0	U	5.0	U	5.0	U	5.0	U	5.0	U
Copper	CU	ICP	IPO92771	25.0	2.0	2.0	U	2.0	U	2.0	U	2.0	U	2.0	U	2.0	U
Lead	PB	PMS	MS100682	3.0	0.1	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
Nickel	NI	ICP	IPO92771	40.0	10.0	10.0	U	10.0	U	10.0	U	10.0	U	10.0	U	10.0	U
Selenium	SE	ICP	IPO92771	5.0	50.0	50.0	U	50.0	U	50.0	U	50.0	U	50.0	U	50.0	U
Thallium	TL	ICP	IPO92771	10.0	50.0	50.0	U	50.0	U	50.0	U	50.0	U	50.0	U	50.0	U
Vanadium	V	ICP	IPO92771	50.0	3.0	3.0	U	3.0	U	3.0	U	3.0	U	3.0	U	3.0	U
Zinc	ZN	ICP	IPO92771	20.0	10.0	10.0	U	10.0	U	10.0	U	10.0	U	10.0	U	10.0	U

Calibration Blanks

CLIENT: AMEC Geomatix

PROJECT: ERP 2011 Shoreline I

SDG: TN19

UNITS: ug/L



ANALYTE	EL	METH	RUN	CRDL	IDL	ICB	C	CCB1	C	CCB2	C	CCB3	C	CCB4	C	CCB5	C
Arsenic	AS	PMS	MS100781	10.0	0.2	0.2	v	0.2	v	0.2	v	0.2	v	0.2	v	0.2	v
Lead	PB	PMS	MS100781	3.0	0.1	0.1	v	0.1	v	0.1	v	0.1	v	0.1	v	0.1	v

ICP Interference Check Sample

CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

SDG: TN19



ICS SOURCE: I.V.

RUNID: IPO92771

INSTRUMENT ID: OPTIMA ICP 2

UNITS: ug/L

ANALYTE	ICSA TV	ICSAB TV	ICSAI	ICSAI	%R	ICSAI	ICSA2	ICSA2	ICSA2	ICSA2	%R	ICSA3	ICSA3	%R
Aluminum	2000000	2000000	206000.6	201923.8	101.0	195055.3	196102.2	98.1						
Antimony		1000	37.6	979.0	97.9	26.9	950.8	95.1						
Arsenic		1000	4.3	960.0	96.0	2.6	1015.2	101.5						
Barium		1000	0.8	980.4	98.0	0.5	983.3	98.3						
Beryllium		1000	0.1	960.4	96.0	0.0	1000.1	100.0						
Boron			-6.4	-8.5		-8.0								
Cadmium		1000	1.4	1015.9	101.6	0.8	1013.8	101.4						
Calcium	1000000	1000000	106593.1	103349.4	103.3	101843.4	100570.1	100.6						
Chromium		1000	-0.2	996.6	99.7	-0.1	987.1	98.7						
Cobalt		1000	2.7	995.0	99.5	2.8	984.5	98.5						
Copper		1000	-0.2	1025.2	102.5	-1.2	1009.5	101.0						
Iron	2000000	2000000	201950.0	197833.9	98.9	188682.8	188999.9	94.5						
Lead		1000	-3.7	968.1	96.8	-3.8	1013.7	101.4						
Magnesium	1000000	1000000	106013.5	98317.3	98.3	100062.5	95476.3	95.5						
Manganese		1000	0.4	926.8	92.7	0.2	930.9	93.1						
Molybdenum			6.9	6.1		5.6								
Nickel		1000	-2.3	932.0	93.2	-2.1	961.4	96.1						
Potassium			49.1	222.2		27.5								
Selenium		1000	-38.0	907.0	90.7	-33.6	977.4	97.7						
Silicon			-23.0	-23.4		-19.7								
Silver		1000	-0.6	996.5	99.7	-0.6	1051.3	105.1						
Sodium			16.1	6.2		10.7								
Strontium			10.6	10.1		10.3								
Thallium		1000	13.7	979.1	97.9	14.2	1031.3	103.1						
Tin			-7.3	-7.2		-7.6								
Titanium			1.5	1.5		1.0								
Vanadium		1000	4.3	958.1	95.8	2.4	948.4	94.8						
Zinc		1000	-1.2	930.8	93.1	-0.9	943.5	94.4						

TCP Interference Check Sample

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN19

ICS SOURCE: I.V.

RUNID: MS100682

INSTRUMENT ID: PE ELAN 6000

UNITS: ug/L

ANALYTE	ICSA TV	ICSAB TV	ICSAI	ICSAB1	%R	ICSA2	ICSAB2	%R	ICSA3	ICSAB3	%R
Antimony				0.1	0.1						
Cadmium	20			0.0	20.0	100.0					
Chromium	20			0.8	20.5	102.5					
Cobalt	20			0.1	19.6	98.0					
Copper	20			0.5	19.4	97.0					
Iron	20000	20000	20327.5	20246.7	101.2						
Manganese	20			0.7	20.0	100.0					
Molybdenum	400	400	427.1	421.6	105.4						
Nickel	20			0.7	18.7	93.5					
Silver	20			0.0	19.3	96.5					
Vanadium				0.0	-0.5						
Zinc	20			1.1	19.8	99.0					

TCP Interference Check Sample

CLIENT: AMEC Geomatix
PROJECT: FRP 2011 Shoreline I
SDG: TN19



ICS SOURCE: I.V.
RUNID: MS100781
INSTRUMENT ID: PE ELAN 6000
UNITS: ug/L

ANALYTE	ICSA TV	ICSAB TV	ICSA1	ICSA1	%R	ICSA2	ICSA2	%R	ICSA3	ICSA3	%R
Antimony			0.1	0.1							
Cadmium	20		0.0	0.0		20.0	100.0				
Chromium	20		0.8	0.8		20.2	101.0				
Cobalt	20		0.1	0.1		18.7	93.5				
Copper	20		0.4	0.4		20.2	101.0				
Iron	200000	200000	19767.1	19820.1	99.1						
Manganese	20		0.7	0.7		19.2	96.0				
Molybdenum	400	400	453.7	461.5	115.4						
Nickel	20		0.6	0.6		20.0	100.0				
Selenium			0.0	0.0		0.1					
Silver	20		0.0	0.0		19.3	96.5				
Vanadium			0.0	0.0		-0.4					
Zinc	20		0.9	0.9		20.1	100.5				

**Post Digest Spike
Sample Recovery**

**ANALYTICAL
RESOURCES
INCORPORATED**

CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

ANALYSIS METHOD: ICP

SDG: TN19

UNITS: ug/L

ANALYTE	CLIENT ID	ARI ID	RUNID	SPIKED		SAMPLE RESULT C	Spike Result C	Spike Added	Matrix	%R
				SAMPLE	RESULT					
Selenium	FRP-092011-001A	TN19APOST	IP092771	9962.47	250.00U	10000	Water	99.6		

ICP Serial Dilutions

ANALYTICAL
RESOURCES
INCORPORATED

CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

ANALYSIS METHOD: ICP

SDG: TN19

UNITS: ug/L

ANALYTE	CLIENT ID	ARI ID	MATRIX	RUNID	INITIAL SAMPLE RESULT (I)	SERIAL DILUTION RESULT (S)		% DIFFER- ENCE	Q
						C	C		
Aluminum	FRP-092011-001L	TN19A-L	Water	IP092771	176049.38	170700.45		3.0	
Cadmium	FRP-092011-001L	TN19A-L	Water	IP092771	1.78 u	10.00 u			
Chromium	FRP-092011-001L	TN19A-L	Water	IP092771	170.72	166.70		2.4	
Copper	FRP-092011-001L	TN19A-L	Water	IP092771	272.58	255.35		6.3	
Nickel	FRP-092011-001L	TN19A-L	Water	IP092771	101.26	108.60 B		7.2	
Selenium	FRP-092011-001L	TN19A-L	Water	IP092771	-19.88 u	250.00 u			
Thallium	FRP-092011-001L	TN19A-L	Water	IP092771	12.14 u	250.00 u			
Vanadium	FRP-092011-001L	TN19A-L	Water	IP092771	444.95	437.50		1.7	
Zinc	FRP-092011-001L	TN19A-L	Water	IP092771	360.43	363.10		0.7	

ICP Serial Dilutions

ANALYTICAL
RESOURCES
INCORPORATED

CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

ANALYSIS METHOD: PMS

SDG: TN19

UNITS: ug/L

ANALYTE	CLIENT ID	ARI ID	MATRIX	RUNID	(I)	INITIAL	SERIAL	% DIFFER-	Q
						SAMPLE RESULT	DILUTION RESULT		
Arsenic	FRP-092011-001L	TN19A-L	Water	MS100781	19.20		20.80 E	8.3	
Lead	FRP-092011-001L	TN19A-L	Water	MS100781	15.63		18.90	20.9	

**IDLs and ICP
Linear Ranges**

**ANALYTICAL
RESOURCES
INCORPORATED**

CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

SDG: TN19

UNITS: ug/L

ANALYTE	EL	METH	INSTRUMENT	WAVELENGTH (nm)	GFA BACK- GROUND	CLP CRDL	RL	RL DATE	ICP LINEAR RANGE (ug/L)	ICP LR DATE
Aluminum	AL	ICP	OPTIMA ICP 2	308.22		200	50.0	4/1/2011	250000.0	8/3/2011
Arsenic	AS	PMS	PE ELAN 6000 MS	0.00		10	0.2	4/1/2011		
Cadmium	CD	ICP	OPTIMA ICP 2	228.80		5	2.0	4/1/2011	20000.0	8/3/2011
Chromium	CR	ICP	OPTIMA ICP 2	267.72		10	5.0	4/1/2011	100000.0	8/3/2011
Copper	CU	ICP	OPTIMA ICP 2	324.75		25	2.0	4/1/2011	40000.0	8/3/2011
Lead	PB	PMS	PE ELAN 6000 MS	0.00		3	0.1	4/1/2011		
Nickel	NI	ICP	OPTIMA ICP 2	231.60		40	10.0	4/1/2011	100000.0	8/3/2011
Selenium	SE	ICP	OPTIMA ICP 2	196.02		5	50.0	4/1/2011	20000.0	8/3/2011
Thallium	TL	ICP	OPTIMA ICP 2	190.86		10	50.0	4/1/2011	30000.0	8/3/2011
Vanadium	V	ICP	OPTIMA ICP 2	292.40		50	3.0	4/1/2011	50000.0	8/3/2011
Zinc	ZN	ICP	OPTIMA ICP 2	213.86		20	10.0	4/1/2011	100000.0	8/3/2011

ICP Interelement Correction Factors

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN19



TEC DATE:
9/26/2011

INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	AI	AS	BA	BE	CA	CD	CO	CR	CU	FE
Aluminum	308.22	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Antimony	206.84	0.3618900	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Arsenic	188.98	0.043543	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.8345790	1.1215100	0.0000000	0.0000000
Barium	233.53	0.0000000	0.0000000	0.0000000	0.0000000	-0.0061516	0.0000000	-0.1892080	0.0000000	0.0000000	0.0496666
Beryllium	313.04	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Cadmium	228.80	0.0000000	5.2418600	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.1195910	0.0000000	0.0000000
Calcium	317.93	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.5252460	0.0000000	0.0000000
Chromium	267.72	0.0000000	0.0000000	0.0000000	0.0000000	0.0220343	0.0000000	0.0000000	0.0000000	0.0000000	-0.0461558
Cobalt	228.62	0.0000000	0.0000000	0.0000000	0.1238430	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Copper	324.75	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.2279050	-0.0318969	0.0000000	-0.0514304
Iron	273.96	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-1.6163900	0.0000000	0.0000000
Lead	220.35	-0.2408020	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-2.3072100	1.2452600	0.0597200
Magnesium	279.08	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-1.6380600	-1.2519300	0.0000000
Manganese	257.61	0.0065976	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.0045330
Molybdenum	202.03	0.0000000	0.0000000	0.0000000	0.0000000	0.0103467	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Nickel	231.60	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Potassium	766.49	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Selenium	196.03	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.3514040	0.0000000	0.0000000	0.0000000
Silicon	288.16	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-3.4885200	0.0000000	0.0000000	0.0000000	0.0000000
Silver	328.07	0.0000000	0.0000000	0.0000000	0.0000000	0.0118175	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Sodium	589.59	0.0000000	0.0000000	0.0000000	0.0000000	8.8385500	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Thallium	190.80	0.1350590	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	1.7836600	0.3510820	0.0000000	-0.1171730
Tin	189.93	0.0000000	0.0000000	0.0000000	0.0000000	-0.0244672	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Titanium	334.90	0.0000000	0.0000000	0.0000000	0.0593584	0.0000000	0.0000000	0.1632010	0.0000000	0.0000000	0.0000000
Vanadium	292.40	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-3.9361800	0.0000000	0.0000000	0.0934302
Zinc	206.20	0.0000000	0.0000000	0.0135250	0.0000000	0.0000000	-0.1475900	0.0000000	0.0000000	0.0000000	0.0000000

FORM XI

ICP Interelement Correction Factors

CLIENT: AMEC Geomatrix
 PROJECT: FRP 2011 Shoreline I
 SDG: TN19



IEC DATE: 9/26/2011
 INSTRUMENT ID: OPTIMA ICP 2

ANALYTE	WAVELENGTH	MG	MN	MO	NI	PR	SB	TI	TL	V	ZN
Aluminum	308.22	0.0000000	0.0000000	15.3131000	0.0000000	0.0000000	1.5167500	0.0000000	17.6996000	0.0000000	
Antimony	206.84	0.0000000	0.0000000	0.0000000	-0.4730780	0.0000000	-0.8897510	0.0000000	-3.3546800	0.0000000	
Arsenic	188.98	0.0000000	0.0000000	2.3330800	0.0000000	0.0000000	-5.4412000	0.0000000	0.0000000	0.0000000	
Barium	233.53	0.0000000	0.0000000	0.0000000	0.0766262	0.0000000	0.0000000	0.0000000	0.6419380	0.0000000	
Beryllium	313.04	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.2960780	0.0000000	
Cadmium	228.80	0.0000000	0.0000000	0.0000000	-0.7324130	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	
Calcium	317.93	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	
Chromium	267.72	0.0464480	0.0000000	0.1395070	0.0000000	0.0000000	0.0000000	0.0000000	0.2773470	0.0000000	
Cobalt	228.62	0.0000000	0.0000000	-0.1579570	0.1588330	0.0000000	0.0000000	1.8115900	0.0000000	0.0000000	
Copper	324.75	0.0000000	0.0000000	0.2688440	0.0000000	0.0000000	0.2461180	0.0000000	0.0000000	0.0000000	
Iron	273.96	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	8.4403600	0.0000000	
Lead	220.35	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	
Magnesium	279.08	0.0000000	0.0000000	-4.6256200	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	
Manganese	257.61	0.0039764	0.0000000	0.0000000	-0.2173850	0.0000000	0.0000000	0.0000000	-0.0271775	0.0000000	
Molybdenum	202.03	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	
Nickel	231.60	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.7744280	0.0000000	0.0000000	0.0000000	
Potassium	766.49	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	
Selenium	196.03	0.07780238	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	
Silicon	288.16	0.0000000	0.0000000	-2.7358100	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	
Silver	328.07	0.0000000	0.2442620	0.2419260	0.0000000	0.0000000	-0.0470302	0.0000000	-0.2758080	0.0000000	
Sodium	589.59	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	
Thallium	190.80	0.0000000	-1.4179000	1.9562000	0.0000000	0.0000000	0.0000000	0.0000000	1.2892100	0.0000000	
Tin	189.93	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	-0.5848020	-0.3044710	0.0000000	0.0000000	
Titanium	334.90	0.0000000	0.0000000	0.9873960	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	
Vanadium	292.40	0.0000000	-0.1398510	-0.6804250	0.0000000	0.0000000	0.6004670	0.0000000	0.0000000	0.0000000	
Zinc	206.20	0.0000000	0.2377960	0.0000000	-0.0708227	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000	

Preparation Log

ANALYTICAL
RESOURCES
INCORPORATED

CLIENT: AMEC Geomatrix

ANALYSIS METHOD: ICP

PROJECT: FRP 2011 Shoreline I

ARI PREP CODE: TWC

SDG: TN19

PREPDATE: 9/22/2011

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
FRP-092011-001	TN19A	0.000	50.0	50.0
FRP-092011-001D	TN19ADUP	0.000	50.0	50.0
FRP-092011-001S	TN19ASPK	0.000	50.0	50.0
FRP-092011-002	TN19B	0.000	50.0	50.0
FRP-092011-003	TN19C	0.000	50.0	50.0
FRP-092011-004	TN19D	0.000	50.0	50.0
FRP-092011-005	TN19E	0.000	50.0	50.0
PBW	TN19MB1	0.000	50.0	50.0
LCSW	TN19MB1SPK	0.000	50.0	50.0

Preparation Log



CLIENT: AMEC Geomatrix

ANALYSIS METHOD: PMS

PROJECT: FRP 2011 Shoreline I

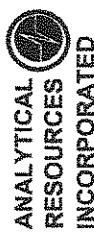
ARI PREP CODE: REN

SDG: TN19

PREPDATE: 9/22/2011

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
FRP-092011-001	TN19A	0.000	50.0	25.0
FRP-092011-001D	TN19ADUP	0.000	50.0	25.0
FRP-092011-001S	TN19ASPK	0.000	50.0	25.0
FRP-092011-002	TN19B	0.000	50.0	25.0
FRP-092011-003	TN19C	0.000	50.0	25.0
FRP-092011-004	TN19D	0.000	50.0	25.0
FRP-092011-005	TN19E	0.000	50.0	25.0
PBW	TN19MB1	0.000	50.0	25.0
LCSW	TN19MB1SPK	0.000	50.0	25.0

Analysis Run Log



CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN19

INSTRUMENT ID: OPTIMA ICP 2

RUNID: IP092771

START DATE: 9/27/2011

END DATE: 9/27/2011

CLIENT ID	ARI ID	DIL.	TIME	%R	AG	AL	B	HA	BE	CA	CD	CO	CRC	CU	FE	HG	K	MG	MN	MO	NA	NI	PB	SB	SE	SI	SN	TI	TL	U	V	ZN
S0	S0		1.00 10250																													
S2	S2		1.00 10291																													
S3	S3		1.00 10305																													
S4	S4		1.00 10332																													
S5	S5		1.00 10353																													
ZZZZZZ	ZZZZZZ		1.00 10400																													
S2	S2		1.00 10435																													
ICV	ICV		1.00 10481																													
ICB	ICB		1.00 10493																													
CRI	CRII		1.00 10531																													
ICSA	ICSAI		1.00 10572																													
ICSAB	ICSABT		1.00 11014																													
CCV	CCV1		1.00 11055																													
CCB	CCB1		1.00 11093																													
PBW	TN19MB1		1.00 11132																													
ZZZZZZ	ZZZZZZ		1.00 11173																													
FRP-092011-001L	TN19A-L		10.00 11214																													
FRP-092011-001	TN19A		25.00 11335																													
FRP-092011-001D	TN19ADUP		5.00 11380																													
FRP-092011-001S	TN19ASEK		5.00 11422																													
FRP-092011-002	TN19B		1.00 11463																													
FRP-092011-003	TN19C		1.00 11503																													
LCSW	TN19MB1SPK		1.00 11550																													
ZZZZZZ	TN16MB2SPK		1.00 11593																													
CCV	CCV2		1.00 12032																													
CCB	CCB2		1.00 12072																													
ZZZZZZ	TN42MB1		1.00 12111																													
ZZZZZZ	TN42A-L		5.00 12152																													
ZZZZZZ	TN42A		1.00 12193																													
ZZZZZZ	TN42ADDP		1.00 12240																													
ZZZZZZ	TN42ASEK		1.00 12282																													
ZZZZZZ	TN42POST		1.00 12325																													
ZZZZZZ	TN42B		1.00 12370																													
ZZZZZZ	TN19D		1.00 12412																													
FRP-092011-004	TN19D		1.00 12453																													

Analysis Run Log



CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN19

INSTRUMENT ID: OPTIMA ICP 2
RUNID: IP092771

START DATE: 9/27/2011
END DATE: 9/27/2011

CLIENT ID	ARI ID	DIL.	TIME	S/R	AG	AL	AS	B	BA	BE	CA	CD	CO	CR	CU	EE	HG	K	MG	MN	MO	NA	NI	PB	SB	SE	SI	TI	TL	U	V	ZN
FRP-092011-005	TN19E	1.00	12495	X																												
ZZZZZZZ	TN42MB1SPK	1.00	12540																													
CCV	CCV3	1.00	12580																													
CCB	CCB3	1.00	13014																													
S0	S0	1.00	13055																													
CCV	CCV4	1.00	13094																													
CCB	CCB4	1.00	13133																													
ZZZZZZZ	TN78MB	1.00	13174																													
ZZZZZZZ	TN78A	1.00	13250																													
ZZZZZZZ	TN79A	1.00	13291																													
ZZZZZZZ	TN42C	1.00	13332																													
ZZZZZZZ	TN42D	1.00	13380																													
ZZZZZZZ	TN42E	1.00	13423																													
ZZZZZZZ	TN42F	1.00	13470																													
ZZZZZZZ	TN42G	5.00	13511																													
ZZZZZZZ	TN42B	2.00	13552																													
ZZZZZZZ	TN78MB3SPK	1.00	13593																													
CCV	CCV5	1.00	14033																													
CCB	CCB5	1.00	14071																													
S2	S2	1.00	14161																													
S3	S3	1.00	14173																													
CCV	CCV6	1.00	14193																													
CCB	CCB6	1.00	14232																													
ZZZZZZZ	TN42C	1.00	14273																													
ZZZZZZZ	TN42D	1.00	14321																													
ZZZZZZZ	TN42E	1.00	14364																													
ZZZZZZZ	TN42F	1.00	14411																													
FRP-092011-001A	TN19APPST	5.00	14452																													
ZZZZZZZ	TN42B	2.00	14493																													
CCV	CCV7	1.00	14534																													
CCB	CCB7	1.00	14573																													
CRI	CRI	1.00	15014																													
ICSA	ICSAF	1.00	15055																													
ICSA	ICSAFB	1.00	15106																													
ICSA	CCV8	1.00	15135																													

Analysis Run Log



CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

SDG: TN19

CLIENT ID	ARI ID	DIL.	TIME	%R	AG AL AS	B	BA	BE	CA	CD	CO	CR	CU	FE	HG	K	MG	MN	MO	NA	NI	PB	SE	SI	SN	TI	TL	U	V	ZN		
CCB	CCB8	1.00	15173																													

INSTRUMENT ID: OPTIMA ICP 2

RUNID: TP092771

METHOD: ICP

START DATE: 9/27/2011

END DATE: 9/27/2011

Analysis Run Log

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN19

INSTRUMENT ID: PE ELAN 6000 MS
RUNID: MS100682 METHOD: PMS

START DATE: 10/6/2011
END DATE: 10/6/2011



CLIENT ID	ARI ID	DIL. TIME	SR	AG AL AS B EA BE CA CD CO CR CU WE HG K MG MN MO NA NI PB SB SE SI SN TI TL U V ZN
S0	S0	1.00 11460		X
ZZZZZZ	ZZZZZZ	1.00 11520		
ZZZZZZ	ZZZZZZ	1.00 11590		X
S3	S3	1.00 12050	X	X
S4	S4	1.00 12120	X	X
S1	S1	1.00 12190	X	X
S2	S2	1.00 12250	X	X
ICV	ICV	1.00 12320	X	X
ICB	ICB	1.00 12400	X	X
CCV	MCCV1	1.00 12450	X	X
CCB	CCB1	1.00 12520	X	X
CRI	MCRI	1.00 12580	X	X
ICSA	ICSAI	1.00 13030	X	X
ICSAB	ICSABI	1.00 13100	X	X
ZZZZZZ	LR200	1.00 13160		
ZZZZZZ	LR300	1.00 13230		
CCV	MCCV2	1.00 13310	X	X
CCB	CCB2	1.00 13380		
ZZZZZZ	TM71MB	2.00 13460		
ZZZZZZ	TM67MBSPK	2.00 13520		
ZZZZZZ	TM71R	2.00 13580		
ZZZZZZ	TM71S	2.00 14050		
ZZZZZZ	TM79C	2.00 14110		
ZZZZZZ	TM67C	2.00 14180		
ZZZZZZ	TM67A	5.00 14240		
ZZZZZZ	TM67B	1000.00 14310		
ZZZZZZ	TM79H	20.00 14370		
ZZZZZZ	TM79J	5.00 14430		
ZZZZZZ	ZZZZZZ			
CCV	MCCV3	1.00 14560	X	X
CCB	CCB3	1.00 15030		
ZZZZZZ	TM79FDUP	1000.00 15150		
ZZZZZZ	TM79F	1000.00 15210		
ZZZZZZ	TM79FSPK	1000.00 15270		
ZZZZZZ	TM79K	1000.00 15330		

Analysis Run Log

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

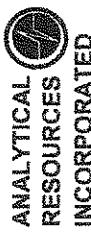
SDG: TN19

INSTRUMENT ID: PE ELAN 6000 MS
RUNID: MS100682 METHOD: PMS

START DATE: 10/6/2011
END DATE: 10/6/2011

CLIENT ID	ARI ID	DIL.	TIME	%R	AG	AL	A8	B	BA	BE	CA	CD	CO	CR	CU	EB	HG	K	MG	NN	MO	NA	NI	PB	SB	SE	SI	SN	T1	T1L	U	V	ZN
ZZZZZZ	TM79G	1000.00	15390																														
ZZZZZZ	TM79G	2.00	15460																														
ZZZZZZ	TN00A-L	100.00	15520																														
ZZZZZZ	TN00A	20.00	15590																														
ZZZZZZ	TN00ADUP	20.00	16050																														
ZZZZZZ	TN00ASEK	20.00	16120																														
CCV	MCCV4	1.00	16220																														
CCB	CCB4	1.00	16300																														
ZZZZZZ	TN00MB2	2.00	16330																														
ZZZZZZ	TN00MB1	20.00	16410																														
ZZZZZZ	TN00MB1SPK	20.00	16460																														
ZZZZZZ	TN00MB2SPK	2.00	16530																														
ZZZZZZ	TN00B-L	50.00	16590																														
ZZZZZZ	TN00B	10.00	17060																														
ZZZZZZ	TN00BDUP	10.00	17120																														
ZZZZZZ	TN00ESK	10.00	17180																														
ZZZZZZ	TN00C	20.00	17250																														
ZZZZZZ	TN00D	5.00	17310																														
CCV	MCCV5	1.00	17370																														
CCB	CCB5	1.00	17440																														
PBW	TN19ME1	2.00	17500																														
LCSW	TN19MB1SPK	2.00	17550																														
ZZZZZZ	TN00E	20.00	18020																														
ZZZZZZ	TN00F	5.00	18080																														
ZZZZZZ	ZZZZZZ	25.00	18140																														
FRP-092011-001	TN19A	5.00	18210																														
FRP-092011-001D	TN19ADUP	5.00	18270																														
FRP-092011-001S	TN19ASEK	5.00	18310																														
FRP-092011-002	TN19B	5.00	18400																														
FRP-092011-003	TN19C	5.00	18470																														
CCV	MCCV6	1.00	18520																														
CCB	CCB6	1.00	18590																														
ZZZZZZ	TN42MB1	2.00	19050																														
ZZZZZZ	TN42MB1SPK	2.00	19100																														
FRP-092011-004	TN19D	5.00	19170																														

Analysis Run Log



CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN19

INSTRUMENT ID: PE ELAN 6000 MS
RUNID: MS100682 METHOD: PMS

START DATE: 10/6/2011
END DATE: 10/6/2011

CLIENT ID	ARI ID	DIL.	TIME	%R	AG AL AS B BA BE CA CD CO CR CU FE HG K MG MN MO NA NI PB SB SE SI SN TI TL U V ZN
FRP-092011-005	TN19E	5.00	19230		
ZZZZZZZ	TN42A-L	50.00	19300		
ZZZZZZZ	TN42A	10.00	19360		
ZZZZZZZ	TN42ADDUP	10.00	19430		
ZZZZZZZ	TN42ASPK	10.00	19490		
ZZZZZZZ	TN42B	5.00	19550		
ZZZZZZZ	TN42C	5.00	20020		
CCV	MCCV7	1.00	20080	X	
CCB	CCB7	1.00	20150	X	X

Analysis Run Log

CITIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SNC-TN19

INSTRUMENT ID: PE ELAN 6000 MS
RUNID: MS100781 METHOD: PMS

START DATE: 10/7/2011
END DATE: 10/7/2011

Analysis Run Log

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN19

INSTRUMENT ID: PE ELAN 6000 MS
RUNID: MS100781 METHOD: PMS

START DATE: 10/7/2011
END DATE: 10/7/2011

CLIENT ID	ARI ID	DIL.	TIME	BR	AG	AL	AS	S	BA	BE	CA	CD	CO	CR	CU	FE	HG	K	MG	MN	MO	NA	NI	PB	SB	SE	SI	SN	TI	TL	U	V	ZN
ZZZZZZ	TN24MB1SPK	2.00	124.60																														
ZZZZZZ	TN24MB3SPK	2.00	125.20																														
ZZZZZZ	TN24MB2SPK	20.00	125.90																														
ZZZZZZ	TN24C	50.00	130.50																														
ZZZZZZ	TN24D	20.00	131.20																														
CCV	MCCV4	1.00	131.80																														
CCB	CCB4	1.00	132.50																														
	TN19A	10.00	133.30																														
ZZZZZZ	TN65A	2.00	133.90																														
ZZZZZZ	TN65B	2.00	134.40																														
ZZZZZZ	TN24B	20.00	135.10																														
ZZZZZZ	TN24E	100.00	135.70																														
ZZZZZZ	TN24F	20.00	140.40																														
ZZZZZZ	TN24B	2.00	141.00																														
ZZZZZZ	TN24E	2.00	141.60																														
ZZZZZZ	TN24F	2.00	142.30																														
FRP-092011-001L	TN19A-L	50.00	143.00																														
CCV	MCCV5	1.00	143.50																														
CCB	CCB5	1.00	144.20																														

Mercury Analysis
Report and Summary QC Forms

ARI Job ID: TN19, TN21

Cover Page**INORGANIC ANALYSIS DATA PACKAGE****ANALYTICAL
RESOURCES
INCORPORATED**

CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

SDG: TN21

CLIENT ID	ARI ID	ARI LIMS ID	REPREP
FRP-092011-001	TN21A	11-20551	
FRP-092011-001D	TN21ADUP	11-20551	
FRP-092011-001S	TN21ASPK	11-20551	
FRP-092011-002	TN21B	11-20552	
PBW	TN21MB1	11-20552	
LCSW	TN21MB1SPK	11-20552	
FRP-092011-003	TN21C	11-20553	
FRP-092011-004	TN21D	11-20554	
FRP-092011-005	TN21E	11-20555	

Were ICP interelement corrections applied ? Yes/No YES

Were ICP background corrections applied ? Yes/No YES

If yes - were raw data generated before application of background corrections ? Yes/No NO

Comments:

THIS DATA PACKAGE HAS BEEN REVIEWED AND AUTHORIZED FOR RELEASE BY:

Signature:

Name: Jay Kuhn

Date:

Title: Inorganics Director

INORGANICS ANALYSIS DATA SHEET
Total Mercury by Method SW7470A

ANALYTICAL
RESOURCES
INCORPORATED

Data Release Authorized: *[Signature]*
Reported: 09/26/11
Date Received: 09/20/11
Page 1 of 1

QC Report No: TN21-AMEC Geomatrix
Project: FRP 2011 Shoreline Investigation
8769

Client/ ARI ID	Date Sampled	Matrix	Prep Date Anal Date	RL	Result
FRP-092011-001 TN21A 11-20551	09/20/11	Water	09/22/11 09/23/11	400	2,840 J
FRP-092011-002 TN21B 11-20552	09/20/11	Water	09/22/11 09/23/11	20.0	181 J
FRP-092011-003 TN21C 11-20553	09/20/11	Water	09/22/11 09/23/11	20.0	175 J
FRP-092011-004 TN21D 11-20554	09/20/11	Water	09/22/11 09/23/11	20.0	193 J
FRP-092011-005 TN21E 11-20555	09/20/11	Water	09/22/11 09/23/11	20.0	20.0 U
MB-092211 Method Blank	NA	Water	09/22/11 09/23/11	20.0	20.0 U

Reported in ng/L

RL-Analytical reporting limit
U-Undetected at reported detection limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Lab Sample ID: TN21A

LIMS ID: 11-20551

Matrix: Water

Data Release Authorized:

Reported: 09/26/11

Sample ID: FRP-092011-001

MATRIX SPIKE

QC Report No: TN21-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation

8769

Date Sampled: 09/20/11

Date Received: 09/20/11

MATRIX SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Spike	Spike Added	% Recovery	Q
Mercury	7470A	2,840	4,320	2,000	74.0%	N

Reported in ng/L

N-Control Limit Not Met

H-% Recovery Not Applicable, Sample Concentration Too High

NA-Not Applicable, Analyte Not Spiked

Percent Recovery Limits: 75-125%

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: FRP-092011-001

DUPPLICATE

Lab Sample ID: TN21A

LIMS ID: 11-20551

Matrix: Water

Data Release Authorized

Reported: 09/26/11

QC Report No: TN21-AMEC Geomatrix

Project: FRP 2011 Shoreline Investigation

8769

Date Sampled: 09/20/11

Date Received: 09/20/11

MATRIX DUPLICATE QUALITY CONTROL REPORT

Analyte	Analysis Method	Sample	Duplicate	RPD	Control Limit	Q
Mercury	7470A	2,840	3,160	10.7%	+/- 20%	

Reported in ng/L

--Control Limit Not Met

L-RPD Invalid, Limit = Detection Limit

INORGANICS ANALYSIS DATA SHEET

TOTAL METALS

Page 1 of 1

Sample ID: LAB CONTROL

Lab Sample ID: TN21LCS
 LIMS ID: 11-20552
 Matrix: Water
 Data Release Authorized: ✓
 Reported: 09/26/11

QC Report No: TN21-AMEC Geomatrix
 Project: FRP 2011 Shoreline Investigation
 8769
 Date Sampled: NA
 Date Received: NA

BLANK SPIKE QUALITY CONTROL REPORT

Analyte	Analysis Method	Spike Found	Spike Added	% Recovery	Q
Mercury	7470A	237	200	118%	

Reported in ng/L

N-Control limit not met
 Control Limits: 80-120%

Calibration Verification

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN21

UNITS: ng/L

ANALYTE	EL	M	RUN	ICVTV	ICV	%R	CCVTV	CCV1	%R	CCV2	%R	CCV3	%R	CCV4	%R	CCV5	%R
Mercury	HG	CVL	HG092301	500.0	474.00	94.8	500.0	494.00	98.8	492.00	98.4	484.00	96.8				

Control Limits: Mercury 80-120; Other Metals 90-110

FORM II (1)

CRDL Standard

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN21

UNITS: ng/L

ANALYTE	EL	M	RUN	CRA/I TV	CR-1	%R	CR-2	%R	CR-3	%R	CR-4	%R	CR-5	%R	CR-6	%R
Mercury	HG	CVL	HG092301	20.0	15.20	76.0										

Control Limits: no control limits have been established by the EPA at this time.

FORM II (2)

Calibration Blanks

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN21



UNITS: ng/L

ANALYTE	EL	METH	RUN	CRDL	IDL	ICB	C	CCB1	C	CCB2	C	CCB3	C	CCB4	C	CCB5	C
Mercury	HG	CVL	HG092301	25.0	20.0	20.0	V										

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**IDLs and ICP
Linear Ranges**

**ANALYTICAL
RESOURCES** 
INCORPORATED

CLIENT: AMEC Geomatrix

PROJECT: FRP 2011 Shoreline I

SDG: TN21

UNITS: ng/L

ANALYTE	EL	METH	INSTRUMENT	WAVELENGTH (nm)	GFA		RL	RL DATE	ICP LINEAR RANGE (ng/L)	ICP LR DATE
					BACK- GROUND	CLP CRDL				
Mercury	HG	CVL	CETAC MERCURY	253.70		25	20.0	4/1/2011		

Preparation Log

ANALYTICAL
RESOURCES
INCORPORATED

CLIENT: AMEC Geomatrix

ANALYSIS METHOD: CVL

PROJECT: FRP 2011 Shoreline I

ARI PREP CODE: TLM

SDG: TN21

PREPDATE: 9/22/2011

CLIENT ID	ARI ID	MASS (g)	INITIAL VOLUME (mL)	FINAL VOLUME (mL)
FRP-092011-001	TN21A	0.000	1.0	20.0
FRP-092011-001D	TN21ADUP	0.000	1.0	20.0
FRP-092011-001S	TN21ASPK	0.000	1.0	20.0
FRP-092011-002	TN21B	0.000	20.0	20.0
FRP-092011-003	TN21C	0.000	20.0	20.0
FRP-092011-004	TN21D	0.000	20.0	20.0
FRP-092011-005	TN21E	0.000	20.0	20.0
PBW	TN21MB1	0.000	20.0	20.0
LCSW	TN21MB1SPK	0.000	20.0	20.0

Analysis Run Log

CLIENT: AMEC Geomatix

PROJECT: FRP 2011 Shoreline I

SDG: TN21

INSTRUMENT ID: CETAC MERCURY
RUNID: HG092301
METHOD: CVL

START DATE: 9/23/2011
END DATE: 9/23/2011

CLIENT ID	ART ID	BIL.	TIME	%R	AC	AL	AS	B	BA	BE	CA	CD	CO	CR	CU	FN	HG	K	MG	MN	MO	NA	NT	PB	SB	SE	SI	SN	TI	TL	U	V	ZN
SO	SO		1.00 16161																														
S20	S20		1.00 16185																														
S50	S50		1.00 16213																														
S100	S100		1.00 16242																														
S200	S200		1.00 16270																														
S400	S400		1.00 16294																														
S1000	S1000		1.00 16323																														
ICV	ICV		1.00 16365																														
TCB	TCB		1.00 16393																														
CCV	CCV		1.00 16421																														
CCB	CCB		1.00 16450																														
CRA	CRA		1.00 16474																														
ZZZZZZ	TM90MB1		1.00 16502																														
ZZZZZZ	TM90MB1SPK		1.00 16530																														
ZZZZZZ	TM90A		1.00 16554																														
ZZZZZZ	TM90ADUP		1.00 16582																														
ZZZZZZ	TM90ASPK		1.00 17011																														
ZZZZZZ	TM90MB2		1.00 17035																														
ZZZZZZ	TM90MB2SPK		1.00 17063																														
ZZZZZZ	TM90B		1.00 17091																														
ZZZZZZ	TM90BDUP		1.00 17120																														
CCV	ACCV2		1.00 17144																														
CCB	CCB2		1.00 17173																														
ZZZZZZ	TM90BSPK		1.00 17201																														
PBW	TN21MB1		1.00 17225																														
LCSW	TN21MB1SPK		1.00 17253																														
FRP-092011-001	TN21A		1.00 17281																														
FRP-092011-001D	TN21ADUP		1.00 17305																														
FRP-092011-001S	TN21ASPK		1.00 17333																														
FRP-092011-002	TN21B		1.00 17361																														
FRP-092011-003	TN21C		1.00 17390																														
FRP-092011-004	TN21D		1.00 17414																														
FRP-092011-005	TN21E		1.00 17442																														
CCV	ACCV3		1.00 17471																														
CCB	CCB3		1.00 17495																														

**General Chemistry Analysis
Report and Summary QC Forms**

ARI Job ID: TN19, TN21

INORGANICS ANALYSIS DATA SHEET
pH by Method EPA 150.1

**ANALYTICAL
RESOURCES
INCORPORATED**

Data Release Authorized: *[Signature]*
 Reported: 09/21/11
 Date Received: 09/20/11
 Page 1 of 1

QC Report No: TN19-AMEC Geomatrix
 Project: FRP 2011 Shoreline Investigation
 8769

Client/ ARI ID	Date Sampled	Matrix	Analysis Date & Batch	RL	Result
FRP-092011-001 TN19A 11-20545	09/20/11	Water	09/20/11 17:08 092011#1	0.01	8.81
FRP-092011-002 TN19B 11-20546	09/20/11	Water	09/20/11 17:08 092011#1	0.01	7.49
FRP-092011-003 TN19C 11-20547	09/20/11	Water	09/20/11 17:08 092011#1	0.01	7.07
FRP-092011-004 TN19D 11-20548	09/20/11	Water	09/20/11 17:08 092011#1	0.01	6.54
FRP-092011-005 TN19E 11-20549	09/20/11	Water	09/20/11 17:08 092011#1	0.01	6.31

Reported in std units

RL-Analytical reporting limit

U-Undetected at reported detection limit

REPLICATE RESULTS-CONVENTIONALS
TN19-AMEC Geomatrix

ANALYTICAL
RESOURCES
INCORPORATED

Matrix: Water
Data Release Authorized:
Reported: 09/21/11

Project: FRP 2011 Shoreline Investiga
Event: 8769
Date Sampled: 09/20/11
Date Received: 09/20/11

Analyte	Date	Units	Sample	Replicate(s)	RPD/RSD
ARI ID: TN19A Client ID: FRP-092011-001					
pH	09/20/11	std units	8.81	8.81	0.00

pH is evaluated as the Absolute Difference between the values rather than
Relative Percent Difference

LAB CONTROL RESULTS-CONVENTIONALS
TN19-AMEC Geomatrix

ANALYTICAL
RESOURCES
INCORPORATED

Matrix: Water
Data Release Authorized
Reported: 09/21/11

Project: FRP 2011 Shoreline Investiga
Event: 8769
Date Sampled: NA
Date Received: NA

Analyte	Date/Time	Units	LCS	Spike Added	Recovery
pH	09/20/11	std units	6.98	7.00	0.02

pH is evaluated as the Absolute Difference between the values rather than Percent Recovery.